

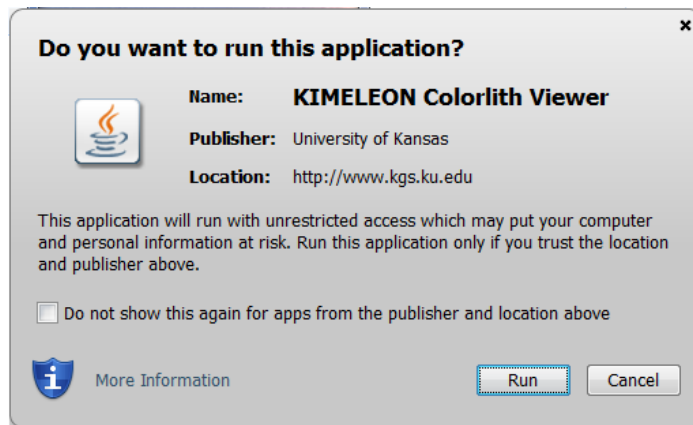
KIMELEON Java Applet

by John R. Victorine

Introduction

The KIMELEON web app has 2 sources for importing well data, 1) the user's PC or 2) the Kansas Geological Survey (KGS) Server & ORACLE Database. This program allows the user to import Log and Tops Data.

There will be occasions when the user would like to inspect the contents of a LAS file. So, for example, the user may wish to find the identity of the well, the types of logs, depth range, well or log parameters. Alternatively, the user can suspect that the file is not coded correctly in LAS format if the Import LAS function fails, and wishes to examine the file to troubleshoot the problem. The LAS file is an ASCII file and can be read by any text editor, i.e. Notepad, WordPad, TextPad, etc. The CSV (Comma Separated Values) files are also ASCII and can be edited in the same way or with Microsoft Excel.



To access Synthetic Seismic web app go to <http://www.kgs.ku.edu/stratigraphic/KIMELEON/>. At the top of the web page there is a menu "Main Page|Description|Applet|Help|Copyright & Disclaimer|". Select the "Applet" menu option a "Warning - Security" Dialog will appear. The program has to be able to read and write to the user's PC and access the Kansas Geological Survey (KGS) Database and File Server, ORACLE requires this dialog. The program does not save your files to KGS, but allows you to access the KGS for well information that may be missing in your Kansas logs. The program does not use Cookies or any hidden software it only reads the LAS and CSV files for the KIMELEON Session and writes Portable Network Graphics (PNG) and/or Portable Document Files (PDF) Files to your. The blue shield on the warning dialog is a symbol that the Java web app is created by a trusted source, which is the University of Kansas.

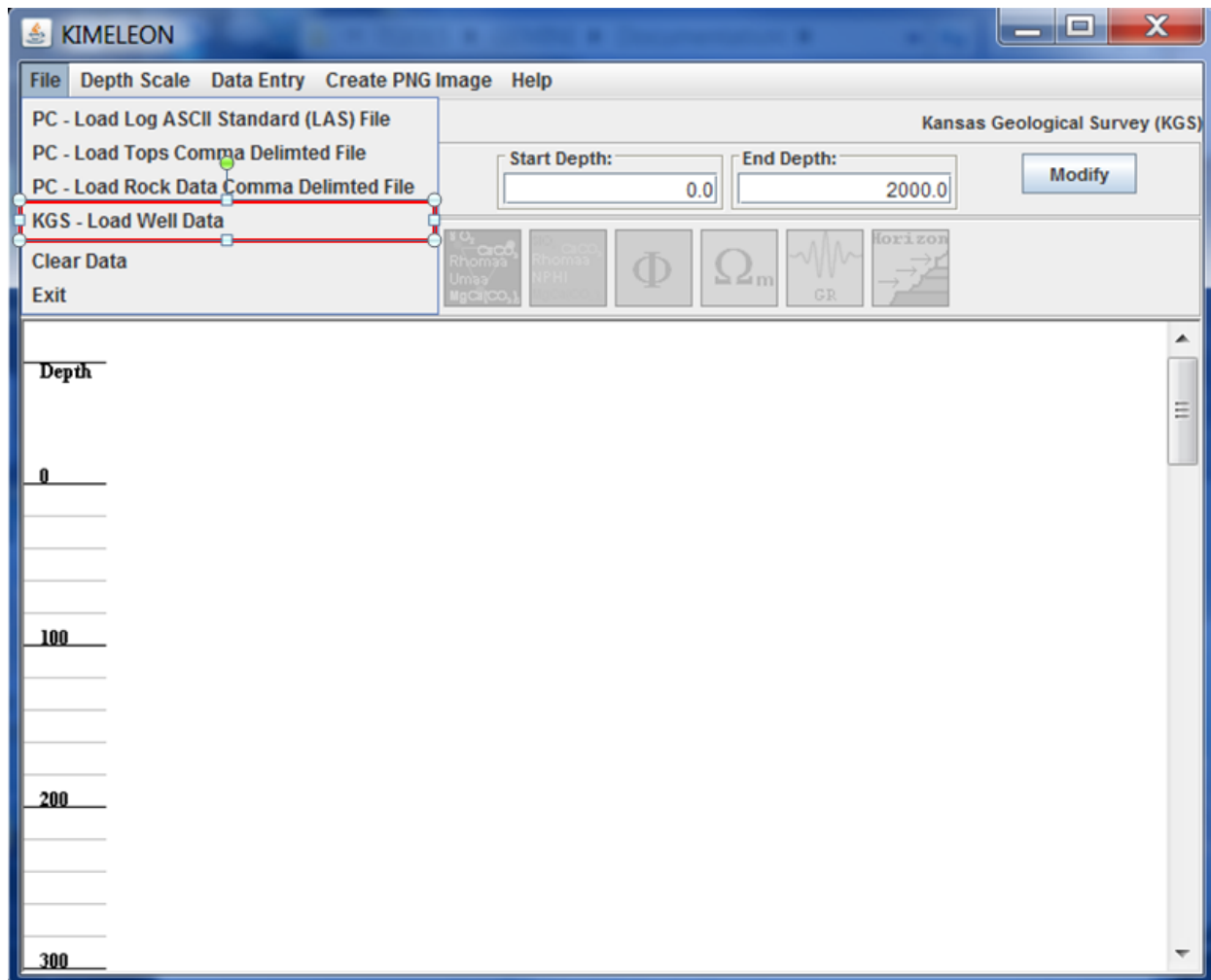
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Importing KGS (Database & Server) Data

KGS (Database & Server) - Importing Well Data

The Kansas Geological Survey (KGS) has a good collection of well data stored in the ORACLE Database and File Server as Files Log ASCII Standard (LAS) version 3.0 Files. In this example the user will download the well data available from the KGS, Log data (LAS version 2.0 File), Tops Data, Measured Core Data, and Perforation Data. The ORACLE Database is accessed by making Stored Procedure PL/SQL calls to the ORACLE Database from which an Extensible Markup Language (XML) data stream is created containing the well data that is passed back to the web app making the request.



Click on the “File” Menu in the menu bar at the top of the “KIMELEON” Dialog, which will display a list of menu options to import well data into the web app. Select the “KGS - Load Well Data” menu option.

Search for Data on KGS Server

Search for Data in Kansas Geological Survey Database:

Search By:

☐ API-Number ☒ Lease name ☐ Township Range Section

Enter Lease (Drop Well Number, Not Case Sensitive):

Newby

Search

List of Oil & Gas Wells:

LTCI	API-Number	Well Name	Operator
LT	15-067-20128	Newby Gas Unit 'C' 2	Amoco Production Co.
T	15-151-21404	Newby 'B' 1-10	Texas Energies, Inc.
T	15-151-20673	Newby 1-14	Texas Energies, Inc.
T	15-189-20067	NEWBY 2-34	MESA PET
T	15-189-20962	NEWBY 3-34	MESA OPERATING
LTCI	15-189-22225	Newby 2-28R	Pioneer Natural Resou
T	15-189-20373	J. T. NEWBY Gas Unit 2	TEXACO INC
	15-125-30191	NEWBY TRUST 'D' 4-6	Dart Cherokee Basin O
	15-205-20826	ORLAND NEWBY 1	M & O OIL COMPANY
	15-205-20894	M & O (MORRIS?) (NEWBY?) 7 TWIN	M & O OIL COMPANY
T	15-155-02712	NEWBY 2	EL DORADO REFINING

Load Data:

LAS File Data Top Picks

Close

Search for Well Data in KGS Database Search By:

• **API-Number** – The user can search the KGS Database for well data by API-Number. The Format for the API is SS-CCC-99999 where

- SS – Two Digit State Code
- CCC – Three Digit County Code
- 99999 – 5 Digit Well Number

• **Lease Name** – The user can search for well data by lease partial phrase, i.e. “Newby”, which will look for all wells with the phrase “Newby” in the lease name.

• **Township-Range-Section** – Search for a list of Wells by a specific area.

List of Kansas wells that match the search criteria

Load Well Data Buttons

- LAS File Data – Load Log ASCII Standard (LAS) Files
- Tops Picks – Load Formation Tops Picks

NOTE: LTCI Column in Table: L-LAS Files; T-Formation Tops; C-Measured Core Data; I-Core Images

This will display the “Search for Data on KGS Server” Dialog, see above image. This dialog allows the user to search the KGS database for well data. In this example, the well of interest will be the Newby 2-28R, this well contains all the well data that can be retrieved from the KGS Database, i.e. Log Data (LAS version 3.0 File), Tops Data, Core Data, and Perforations.

As the Summary image suggests there are 3 methods for searching for the well data within this dialog,

- By API-Number – KGS has a specific format for the API-Number, i.e. SS-CCC-99999 where SS is the state code for Kansas 15, CCC is the county code for Newby 2-28R it is 189 for Stevens County and the 5-Digit Well Number for Newby 2-28R it is 22225.

Search By:

☒ API-Number ☐ Lease name ☐ Township Range Section

Enter API-Number :

15-189-22225

Search

- By Partial Lease Name – The stored procedure used to retrieve the well header information allows the user to enter a partial phrase, in this example Newby. The program places a “%” in front and back of the phrase and sends the request to the Database, i.e. “%Newby%”.

Search By:

☐ API-Number
 ☒ Lease name
 ☐ Township Range Section

Enter Lease (Drop Well Number, Not Case Sensitive):

Newby

Search

- By Township Range Section – This search is by location in Kansas, this search also allows the user to enter just the Township and Range to search for wells, e.g. to look for the Newby 2-28R, enter Township as 31 set the S (South) Radio button and Range as 37 set the W (West) Radio button.

Search By:

☐ API-Number
 ☐ Lease name
 ☒ Township Range Section

Section:
 Township:
☐ N ☒ S
 Range:
☒ W ☐ E

Search

The user only needs to enter the above data and select the “Search” Button to display the list of Wells in the Kansas Database that match the search criteria. In the image below the Lease Name “Newby” was entered to search for all wells in Kansas with the Phrase Newby in it. The user searches through the list until they find the well of interest. In this example it is the Newby 2-28R, which is highlighted.

Search for Data on KGS Server

Search for Data in Kansas Geological Survey Database:

Search By:

☐ API-Number ☒ Lease name ☐ Township Range Section

Enter Lease (Drop Well Number, Not Case Sensitive):

Newby

Search

List of Oil & Gas Wells:

LTCI	API-Number	Well Name	Operator
LT	15-067-20128	Newby Gas Unit 'C' 2	Amoco Production Co.
T	15-151-21404	Newby 'B' 1-10	Texas Energies, Inc.
T	15-151-20673	Newby 1-14	Texas Energies, Inc.
T	15-189-20067	NEWBY 2-34	MESA PET
T	15-189-20962	NEWBY 3-34	MESA OPERATING
LTCI	15-189-22225	Newby 2-28R	Pioneer Natural Resou
T	15-189-20373	J. T. NEWBY Gas Unit 2	TEXACO INC
	15-125-30191	NEWBY TRUST 'D' 4-6	Dart Cherokee Basin O
	15-205-20826	ORLAND NEWBY 1	M & O OIL COMPANY
	15-205-20894	M & O (MORRIS?) (NEWBY?) 7 TWIN	M & O OIL COMPANY
T	15-155-02712	NEWBY 2	EL DORADO REFINING

Load Data:

LAS File Data Top Picks

Close

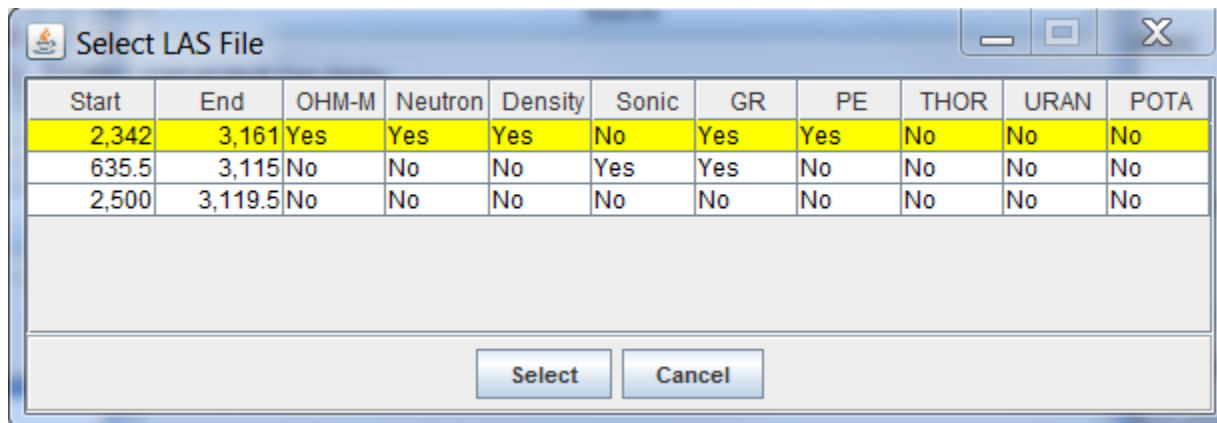
Notice that the LTCI represents the type of data that the well contains. It is a visual aid that lets the user see what is available before trying to download the data. If you require a LAS file you would want to see an L in that column. The LTCI labels stand for the following,

- L – Log ASCII Standard (LAS) version 2.0 Files
- T – Tops Data (Stratigraphic Unit Horizons)
- C – Measured Core Data
- I – Core Joint Photographic Experts Group (JPEG) Image Files

This dialog allows the user to now download each of the data types that are available.

Load KGS Well Data – LAS File Data

The “Search for Data on KGS Server” Dialog allows the user to download data from the KGS Database & Server to the web app. The “LAS File Data” Button will display the “Select LAS File” Dialog with a list of LAS version 2.0 Files that are available.

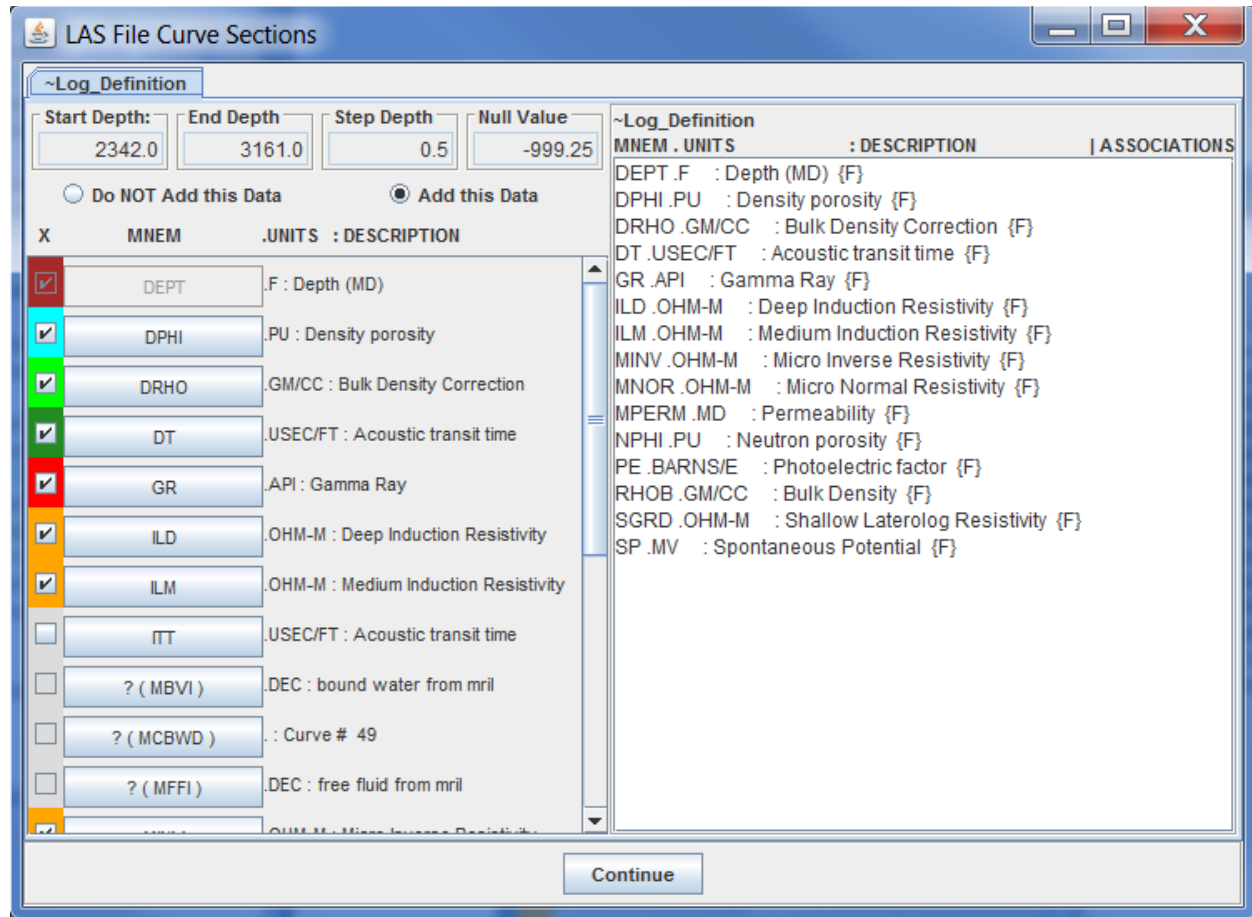


In this example there are three LAS files available, with a table suggesting the log data type in the file. In the beginning of the GEMINI Project (2000-2003) a precursor to the GEMINI Tools the KGS populated the Tool Types from every log that was in the KGS Server at that time. Unfortunately KGS has not maintain that table for wells uploaded after 2003 so the LAS File may have “No” for all the log types, which is not accurate. The user will need to open or download the file or search for the Well on the KGS Master List of Oil and Gas Wells in Kansas Web Page (<http://www.kgs.ku.edu/Magellan/Qualified/index.html>) to see what is in the File Header before deciding to download data from this program. For this example the first log has most of the data necessary except the Spectral Gamma Ray Logs. The Table above identifies the following log types,

- OHM-M – Resistivity Logs
- Neutron – Neutron Porosity Log
- Density – Bulk Density and/or Density Porosity Log
- Sonic – Acoustic Transit Time and/or Sonic Porosity Log
- GR – Gamma Ray (API units) Log
- PE – Photoelectric Factor Log
- THOR – Thorium Concentration
- URAN – Uranium Concentration
- POTA – Potassium Concentration)

In this example the first log contains the data needed, highlight the first log and click on the “Select” Button to display the “LAS File Curve Sections” Dialog. The “LAS File Curve Sections” Dialog allows the user to map unknown LAS Curve Mnemonics to the KGS “Standard” Curve Mnemonics so they will be plotted in the Synthetic Seismic Web App Plot. This program reads the “LAS Tool Curve Mnemonics map to KGS Standard Mnemonics” XML File (http://www.kgs.ku.edu/software/gemini/data/las_standard_tools.xml), which will

automatically maps the Curve Mnemonics from the LAS file to one of 31 KGS “Standard” Curve Mnemonics.



As you can see this log has all the log types of interest, Gamma Ray API, Resistivity, Neutron/Density, Photoelectric Factor, Sonic and Permeability. If a curve Mnemonic is not recognized the program will place a “?” in front of the Mnemonic, e.g. “?(MPPM16)” for the “: Curve # 51” Log Curve. If the user is satisfied with the automatic curve selections, which are checked and color coded, they only need to select the “Continue” Button at the bottom of the Dialog to import the file. The next section will take the user through a series of examples in changing the curve selections and mapping unknown curve mnemonics.

Notice that some of the check boxes are colored with different colors, which shows that the curves were automatically selected, but also to represent the curve type by color. The Curves are colored by type (data units) as follows,

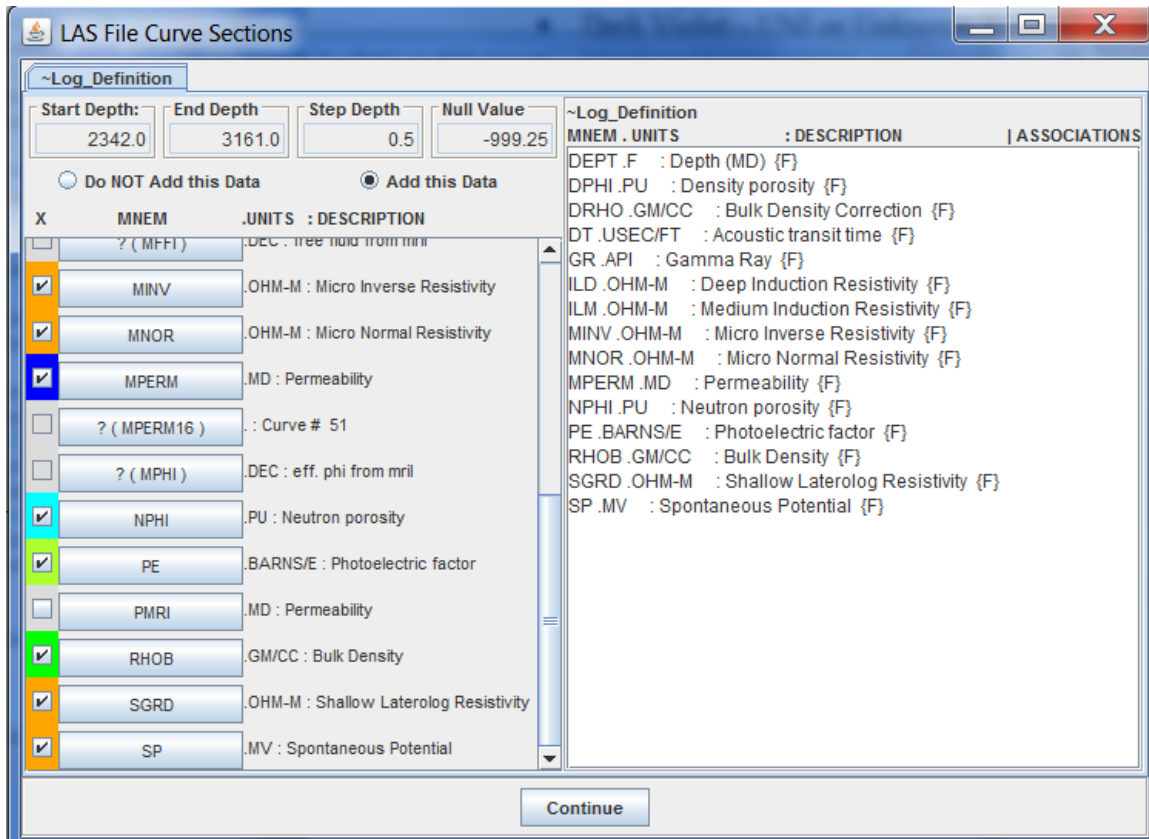
- Orange - OHM-M or Resistivity Logs
- Cyan – PU or porosity Logs, Neutron Porosity, Density Porosity, etc.
- Greenish yellow – BARNS/E or Photoelectric Factor Logs
- Green – GM/CC or Bulk Density Log
- Forest Green – USEC/FT or the Acoustic Transit Time Log

- Red – API, PPM or “%” as Radioactive logs, Gamma Ray, Spectral Gamma Ray, etc.
- Blue – MD or Permeability Logs
- Brown – F, FT or IN or Depth
- Middle yellow – FRAC, or other log curve types.
- Dark Violet – UNI or Unknown Linear Curves
- Medium Violet – UNL or Unknown Logrithum Curves

The color coding of the selected curves were added to also help the user visually recognize that a curve was selected or not.

Map Curves & Change Curve Selections

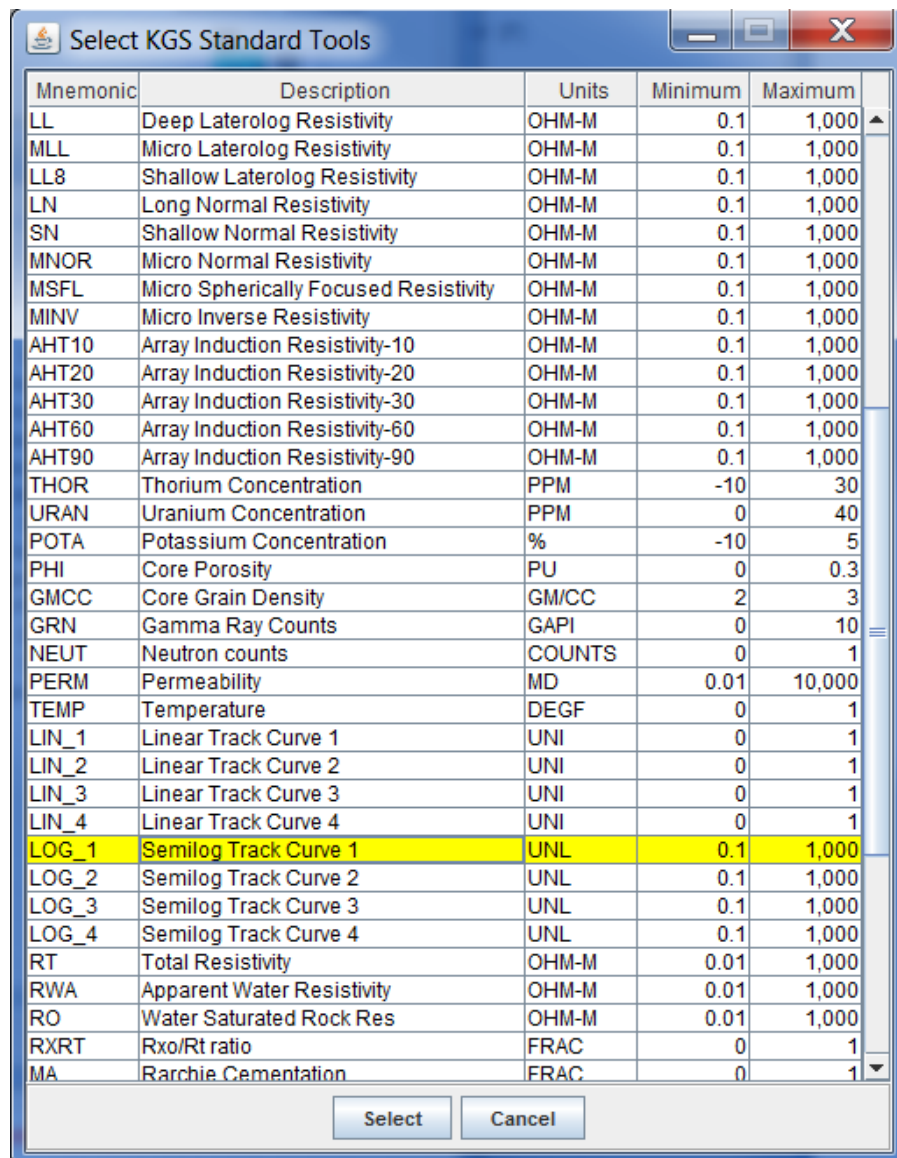
Some logs will have curve mnemonics that are not recognized as one of the KGS “Standard” Curve Mnemonics. The user will need to map the log curve to one of the KGS standard curves if they want to display the curve. Slide the scroll bar down to the Permeability Curves MPERM and ?(MPERM16).



Click on the “?(MPERM16)” Button to display the “Select KGS Standard Tools” Dialog. This dialog provides a list of the KGS “Standard” Curve Mnemonics, from which the user can map an unrecognized log curve to one of the KGS standard curve mnemonics. The KGS “Standard” Curve Mnemonics List was created as a way to standardize the alpha bit soup of Log Mnemonics. Each logging company has their own curve mnemonics to represent similar tools.

The Synthetic Seismic Web App program is a later version of code from the GEMINI Project Synthetic Seismic Web App Module, which needed to standardize the log curves so the curves could be automatically read and assigned a plot track. The “LAS Tool Curve Mnemonics map to KGS Standard Mnemonics” XML File was created to map the log curves from logs that were part of the KGS LAS File Collection which is not a complete list of possible curve mnemonics.

To map the unknown curve mnemonic “?(MPERM16)” you first notice that it is similar to the MPERM curve above, which is a permeability curve. In this example both curves will be plotted together, but the Permeability Plot Track can only plot one curve. The web app has the ability to allow the user to plot up to 4 unknown logarithm curves and 4 unknown linear curves. The permeability is usually plotted as logarithmic. Click on the “?(MPERM16)” Button to display the



“Select KGS Standard Tools” Dialog. Slide the scroll bar down to the “LOG_1” Mnemonic – Semilog Track Curve 1 and highlight that curve. Click on the “Select” Button to map the “?(MPERM16)” to the Semilog Track Curve 1.

LAS File Curve Sections

~Log_Definition

Start Depth: 2342.0 End Depth: 3161.0 Step Depth: 0.5 Null Value: -999.25

☐ Do NOT Add this Data ☒ Add this Data

X	MNEM	.UNITS : DESCRIPTION
<input type="checkbox"/>	? (MFFI)	.DEC : free fluid from mrii
<input checked="" type="checkbox"/>	MINV	.OHM-M : Micro Inverse Resistivity
<input checked="" type="checkbox"/>	MNOR	.OHM-M : Micro Normal Resistivity
<input checked="" type="checkbox"/>	MPERM	.MD : Permeability
<input type="checkbox"/>	MPERM16	.UNL : () Curve # 51
<input type="checkbox"/>	? (MPHI)	.DEC : eff. phi from mrii
<input checked="" type="checkbox"/>	NPHI	.PU : Neutron porosity
<input checked="" type="checkbox"/>	PE	.BARNS/E : Photoelectric factor
<input type="checkbox"/>	PMRI	.MD : Permeability
<input checked="" type="checkbox"/>	RHOB	.GM/CC : Bulk Density
<input checked="" type="checkbox"/>	SGRD	.OHM-M : Shallow Laterolog Resistivity
<input checked="" type="checkbox"/>	SP	.MV : Spontaneous Potential

Continue

~Log_Definition

MNEM . UNITS	: DESCRIPTION	ASSOCIATIONS
DEPT .F	: Depth (MD) {F}	
DPHI .PU	: Density porosity {F}	
DRHO .GM/CC	: Bulk Density Correction {F}	
DT .USEC/FT	: Acoustic transit time {F}	
GR .API	: Gamma Ray {F}	
ILD .OHM-M	: Deep Induction Resistivity {F}	
ILM .OHM-M	: Medium Induction Resistivity {F}	
MINV .OHM-M	: Micro Inverse Resistivity {F}	
MNOR .OHM-M	: Micro Normal Resistivity {F}	
MPERM .MD	: Permeability {F}	
NPHI .PU	: Neutron porosity {F}	
PE .BARNS/E	: Photoelectric factor {F}	
RHOB .GM/CC	: Bulk Density {F}	
SGRD .OHM-M	: Shallow Laterolog Resistivity {F}	
SP .MV	: Spontaneous Potential {F}	

The ?(MPERM16) Curve has been changed to MPERM16 removing the ?() around the Curve Mnemonic. Also select the check box next to it, which changes to a dark violet.

LAS File Curve Sections

~Log_Definition

Start Depth: 2342.0 End Depth: 3161.0 Step Depth: 0.5 Null Value: -999.25

☐ Do NOT Add this Data ☒ Add this Data

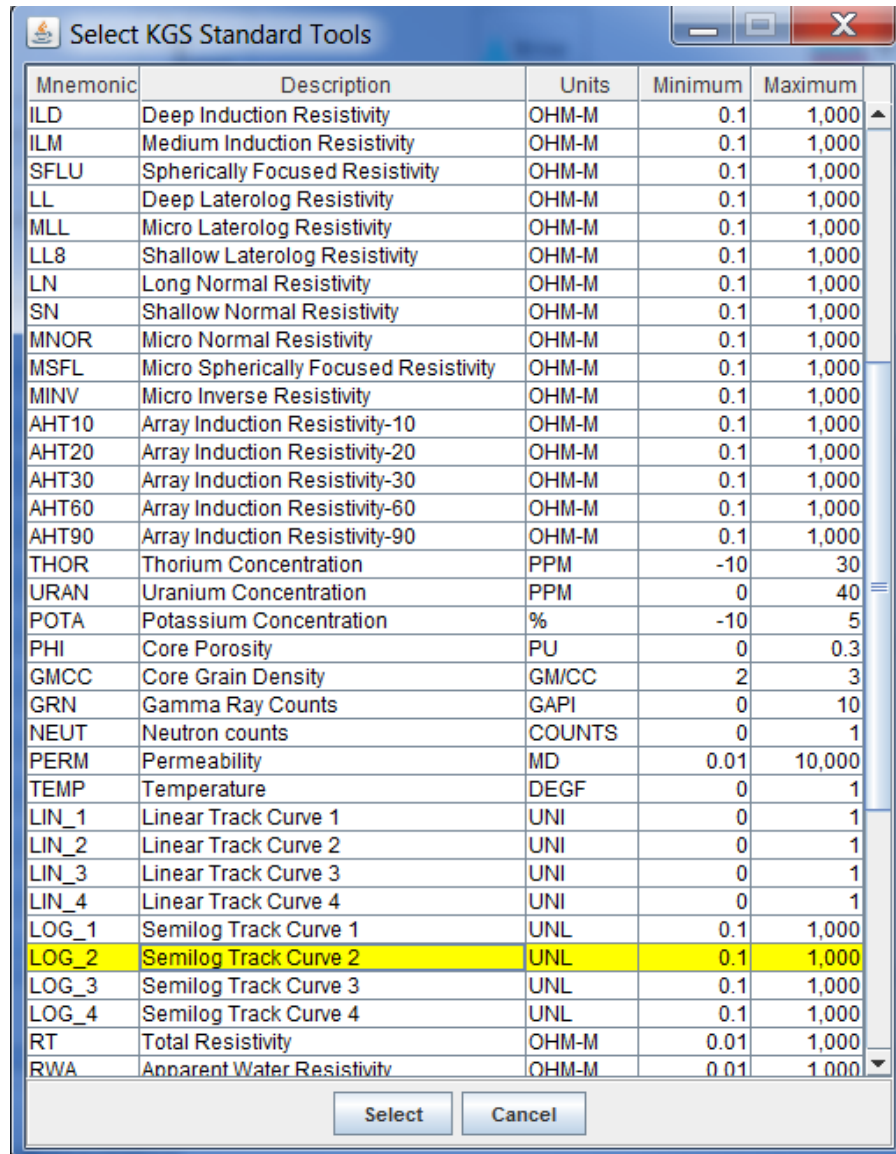
X	MNEM	.UNITS : DESCRIPTION
<input type="checkbox"/>	? (MBVI)	.DEC : bound water from mrii
<input type="checkbox"/>	? (MCBWD)	. : Curve # 49
<input type="checkbox"/>	? (MFFI)	.DEC : free fluid from mrii
<input checked="" type="checkbox"/>	MINV	.OHM-M : Micro Inverse Resistivity
<input checked="" type="checkbox"/>	MNOR	.OHM-M : Micro Normal Resistivity
<input checked="" type="checkbox"/>	MPERM	.MD : Permeability
<input checked="" type="checkbox"/>	MPERM16	.UNL : () Curve # 51
<input type="checkbox"/>	? (MPHI)	.DEC : eff. phi from mrii
<input checked="" type="checkbox"/>	NPHI	.PU : Neutron porosity
<input checked="" type="checkbox"/>	PE	.BARNS/E : Photoelectric factor
<input type="checkbox"/>	PMRI	.MD : Permeability
<input checked="" type="checkbox"/>	RHOB	.GM/CC : Bulk Density

Continue

~Log_Definition

MNEM . UNITS	: DESCRIPTION	ASSOCIATIONS
DEPT .F	: Depth (MD) {F}	
DPHI .PU	: Density porosity {F}	
DRHO .GM/CC	: Bulk Density Correction {F}	
DT .USEC/FT	: Acoustic transit time {F}	
GR .API	: Gamma Ray {F}	
ILD .OHM-M	: Deep Induction Resistivity {F}	
ILM .OHM-M	: Medium Induction Resistivity {F}	
MINV .OHM-M	: Micro Inverse Resistivity {F}	
MNOR .OHM-M	: Micro Normal Resistivity {F}	
MPERM .MD	: Permeability {F}	
MPERM16 .UNL	: Curve # 51 {F}	
NPHI .PU	: Neutron porosity {F}	
PE .BARNS/E	: Photoelectric factor {F}	
RHOB .GM/CC	: Bulk Density {F}	
SGRD .OHM-M	: Shallow Laterolog Resistivity {F}	
SP .MV	: Spontaneous Potential {F}	

In order to plot both Permeability Curves on the same track, the MPERM must be assigned to the unknown Log Curves with MPERM16. Click on the MPERM Mnemonic Button to display the



“Select KGS Standard Tools” Dialog. Slide the scroll bar down to the “LOG_2” Mnemonic – Semilog Track Curve 2 and highlight that curve. Click on the “Select” Button to map the “MPERM to the Semilog Track Curve 2.

LAS File Curve Sections

~Log_Definition

Start Depth: 2342.0 End Depth: 3161.0 Step Depth: 0.5 Null Value: -999.25

☐ Do NOT Add this Data ☒ Add this Data

X	MNEM	.UNITS	DESCRIPTION
<input type="checkbox"/>	? (MBVI)	.DEC	: bound water from mrii
<input type="checkbox"/>	? (MCBWD)	.	: Curve # 49
<input type="checkbox"/>	? (MFFI)	.DEC	: free fluid from mrii
<input checked="" type="checkbox"/>	MINV	.OHM-M	: Micro Inverse Resistivity
<input checked="" type="checkbox"/>	MNOR	.OHM-M	: Micro Normal Resistivity
<input checked="" type="checkbox"/>	MPERM	.UNL : (UNL) (MD)	Permeability
<input checked="" type="checkbox"/>	MPERM16	.UNL : ()	Curve # 51
<input type="checkbox"/>	? (MPHI)	.DEC	: eff. phi from mrii
<input checked="" type="checkbox"/>	NPHI	.PU	: Neutron porosity
<input checked="" type="checkbox"/>	PE	.BARNS/E	: Photoelectric factor
<input type="checkbox"/>	PMRI	.MD	: Permeability

~Log_Definition

MNEM	.UNITS	DESCRIPTION	ASSOCIATIONS
DEPT	.F	: Depth (MD) {F}	
DPHI	.PU	: Density porosity {F}	
DRHO	.GM/CC	: Bulk Density Correction {F}	
DT	.USEC/FT	: Acoustic transit time {F}	
GR	.API	: Gamma Ray {F}	
ILD	.OHM-M	: Deep Induction Resistivity {F}	
ILM	.OHM-M	: Medium Induction Resistivity {F}	
MINV	.OHM-M	: Micro Inverse Resistivity {F}	
MNOR	.OHM-M	: Micro Normal Resistivity {F}	
MPERM	.UNL	: Permeability {F}	
MPERM16	.UNL	: Curve # 51 {F}	
NPHI	.PU	: Neutron porosity {F}	
PE	.BARNS/E	: Photoelectric factor {F}	
RHOB	.GM/CC	: Bulk Density {F}	
SGRD	.OHM-M	: Shallow Laterolog Resistivity {F}	
SP	.MV	: Spontaneous Potential {F}	

Continue

Click the check box next to the MPERM Mnemonic Button you will notice that it has the same color as MPERM16, which now groups the two curves together. Select the “Continue” Button to load the LAS File into the web app.

Load KGS Well Data – Top Picks

The “Search for Data on KGS Server” Dialog allows the user to download data from the KGS Database & Server to the web app. The “Top Picks” Button will display the “Move/Merge KGS Data” Dialog with available top picks grouped by the sources creating the tops.

Move/Merge KGS Data.

KGS Stratigraphic Units:

- ☒ HUG ELOG-EM
- ☐ MKD
- ☐ MKD-07/2006

List of Sources for the tops, e.g. Newby 2-28R has 3 sources of tops data. The user can search through and add some or all to the web app.

Add to User's Stratigraphic Units List:

☐ Remove & Replace ☒ Add to List ☐ Add New Units Only

Source	Top	Base	Name	R
HUG ELOG-EM	0	0	Council Grove	GROL
HUG ELOG-EM	728	0	Day Creek Dolomite	FORM
HUG ELOG-EM	1,090	1,170	Blaine	FORM
HUG ELOG-EM	1,250	1,412	Cedar Hills Sandstone	FORM
HUG ELOG-EM	1,690	1,759	Stone Corral	FORM
HUG ELOG-EM	2,182	2,516	Wellington	FORM
HUG ELOG-EM	2,291	0	Hutchinson Salt	MEMB
HUG ELOG-EM	2,496	0	Hollenberg Limestone	MEMB
HUG ELOG-EM	2,516	2,832	Chase	GROL
HUG ELOG-EM	2,516	2,536	Herington Limestone	MEMB

User's Stratigraphic Units:

Source	Top	Base	Name	Rank	P
HUG ELOG-EM	728	0	Day Creek Dolomite	FORMATION	P
HUG ELOG-EM	1,090	1,170	Blaine	FORMATION	P
HUG ELOG-EM	1,250	1,412	Cedar Hills Sandstone	FORMATION	P
HUG ELOG-EM	1,690	1,759	Stone Corral	FORMATION	P
HUG ELOG-EM	2,182	2,516	Wellington	FORMATION	P
HUG ELOG-EM	2,291	0	Hutchinson Salt	MEMBER	P
HUG ELOG-EM	2,496	0	Hollenberg Limestone	MEMBER	P
HUG ELOG-EM	2,516	2,832	Chase	GROUP	P
HUG ELOG-EM	2,516	2,536	Herington Limestone	MEMBER	P
HUG ELOG-EM	2,536	2,538	Paddock Shale	MEMBER	P
HUG ELOG-EM	2,544	2,580	Krider Limestone	MEMBER	P
HUG ELOG-EM	2,580	2,594	Odell Shale	FORMATION	P
HUG ELOG-EM	2,594	2,629	Winfield Limestone	FORMATION	P
HUG ELOG-EM	2,632	0	Gage Shale	MEMBER	P
HUG ELOG-EM	2,655	2,704	Towanda Limestone	MEMBER	P
HUG ELOG-EM	2,742	2,756	East Diley Limestone	MEMBER	P

Buttons: Add, Add All, Clear Selection, Clear Selection, Remove, Remove All, Load Data, Close.

“Add to User's Stratigraphic Units List” Table shows the tops selected by the source, e.g. “HUG ELOG-EM” Source Tops List.

Radio Buttons

Remove & Replace – move the selected tops and replace any duplicate names

Add to List – move the selected tops to the “User's Stratigraphic Units” Table

Add New Units Only – move on the selected tops that are not already in the “User's Stratigraphic Units” Table

Table Buttons

Add – add the highlighted top(s) to the “User's Stratigraphic Units” Table. Note: this table will allow the user to select multiple wells by using the “Ctrl” Key and the left click of mouse.

Add All – copy the list of tops to the “User's Stratigraphic Units” Table.

Clear Selection – remove the highlight on tops selected.

“Add to User's Stratigraphic Units List” Table.

“User's Stratigraphic Units” Table shows the list of tops that will appear in the web app when the user selects the “Load Data” Button.

Table Buttons

Clear Selection – remove the highlight on tops selected.

Remove – remove the highlighted top(s) from the table. Note: this table will allow the user to select multiple wells by using the “Ctrl” Key and the left click of mouse.

Remove All – remove all tops from the table.

Load Data – transfer the tops list to the web app calling.

Close – Close this dialog

This dialog allows the user to add all or some the tops from each of the sources. Both tables are set up so the user can use the “Ctrl” Key with the left click of mouse to select multiple tops, i.e.

Notice that the only some of the tops are selected. You can then select the Add Button to move only the selected tops to the “User’s Stratigraphic Units” Table.

MKD Source Example:

KGS Stratigraphic Units:

☐ HUG ELOG-EM

☒ MKD

☐ MKD-07/2006

Add to User's Stratigraphic Units List:

☒ Remove & Replace ☐ Add to List ☐ Add New Units Only

Source	Top	Base	Name	R
MKD	0	2,773	Fort Riley Limestone	MEMB
MKD	0	2,693	Towanda Limestone	MEMB
MKD	2,538	2,580	Krider Limestone	MEMB
MKD	2,629	0	Gage Shale	MEMB
MKD	2,712	0	Fort Riley Limestone	MEMB
MKD	2,777	2,789	Florence Limestone	MEMB
MKD	2,807	0	Wreford Limestone	FORM
MKD	2,832	0	Council Grove	GROL
MKD	2,832	0	Council Grove	GROL
MKD	2,832	2,853.5	Speiser Shale	FORM

(1) Select the MKD Source, which will be displayed in the “Add to User’s Stratigraphic Units List” Table.

Add Add All Clear Selection

KGS Stratigraphic Units:

☐ HUG ELOG-EM

☒ MKD

☐ MKD-07/2006

Add to User's Stratigraphic Units List:

☐ Remove & Replace ☒ Add to List ☐ Add New Units Only

Source	Top	Base	Name	R
MKD	0	2,773	Fort Riley Limestone	MEMB
MKD	0	2,693	Towanda Limestone	MEMB
MKD	2,538	2,580	Krider Limestone	MEMB
MKD	2,629	0	Gage Shale	MEMB
MKD	2,712	0	Fort Riley Limestone	MEMB
MKD	2,777	2,789	Florence Limestone	MEMB
MKD	2,807	0	Wreford Limestone	FORM
MKD	2,832	0	Council Grove	GROL
MKD	2,832	0	Council Grove	GROL
MKD	2,832	2,853.5	Speiser Shale	FORM

(2) Select the “Add to List” Radio button.

Add Add All Clear Selection

KGS Stratigraphic Units:

☐ HUG ELOG-EM
☒ MKD
☐ MKD-07/2006

Add to User's Stratigraphic Units List:

☐ Remove & Replace ☒ Add to List ☐ Add New Units Only

Source	Top	Base	Name	R
MKD	0	2,773	Fort Riley Limestone	MEMB
MKD	0	2,693	Towanda Limestone	MEMB
MKD	2,538	2,580	Krider Limestone	MEMB
MKD	2,629	0	Gage Shale	MEMB
MKD	2,712	0	Fort Riley Limestone	MEMB
MKD	2,777	2,789	Florence Limestone	MEMB
MKD	2,807	0	Wreford Limestone	FORM
MKD	2,832	0	Council Grove	GROU
MKD	2,832	0	Council Grove	GROU
MKD	2,832	2,853.5	Speiser Shale	FORM

User's Stratigraphic Units:

Source	Top	Base	Name	Rank	
MKD	2,538	2,580	Krider Limestone	MEMBER	P
MKD	2,629	0	Gage Shale	MEMBER	P
MKD	2,712	0	Fort Riley Limestone	MEMBER	P
MKD	2,777	2,789	Florence Limestone	MEMBER	P
MKD	2,807	0	Wreford Limestone	FORMATION	P
MKD	2,832	0	Council Grove	GROUP	P
MKD	2,832	0	Council Grove	GROUP	P
MKD	2,832	2,853.5	Speiser Shale	FORMATION	P
MKD	2,853.5	2,894.5	Funston Limestone	FORMATION	P
MKD	2,894.5	2,910.5	Blue Rapids Shale	FORMATION	P
MKD	2,910.5	2,929	Crouse Limestone	FORMATION	P
MKD	2,929	2,933.5	Easley Creek Shale	FORMATION	P
MKD	2,933.5	2,947	Middleburg Limestone	MEMBER	P
MKD	2,947	2,957.5	Hooser Shale	MEMBER	P
MKD	2,957.5	2,962.5	Eiss Limestone	MEMBER	P
MKD	2,962.5	2,973.5	Stages Shale	FORMATION	P

Once the list of tops are in the "User's Stratigraphic Units" Table the user can edit the list by removing any duplicate or invalid tops. Notice that the "Council Grove" Top occurs 2 times in the list. Highlight the one of the "Council Grove" tops.

User's Stratigraphic Units:

Source	Top	Base	Name	Rank	
MKD	2,538	2,580	Krider Limestone	MEMBER	P ▲
MKD	2,629	0	Gage Shale	MEMBER	P
MKD	2,712	0	Fort Riley Limestone	MEMBER	P
MKD	2,777	2,789	Florence Limestone	MEMBER	P
MKD	2,807	0	Wreford Limestone	FORMATION	P
MKD	2,832	0	Council Grove	GROUP	P
MKD	2,832	0	Council Grove	GROUP	P
MKD	2,832	2,853.5	Speiser Shale	FORMATION	P
MKD	2,853.5	2,894.5	Funston Limestone	FORMATION	P
MKD	2,894.5	2,910.5	Blue Rapids Shale	FORMATION	P
MKD	2,910.5	2,929	Crouse Limestone	FORMATION	P
MKD	2,929	2,933.5	Easily Creek Shale	FORMATION	P
MKD	2,933.5	2,947	Middleburg Limestone	MEMBER	P
MKD	2,947	2,957.5	Hooser Shale	MEMBER	P
MKD	2,957.5	2,962.5	Eiss Limestone	MEMBER	P
MKD	2,962.5	2,972.5	Stearns Shale	FORMATION	P ▼

Buttons: Clear Selection, Remove, Remove All, Load Data, Close

Now select the "Remove" Button.

User's Stratigraphic Units:

Source	Top	Base	Name	Rank	
MKD	2,538	2,580	Krider Limestone	MEMBER	P ▲
MKD	2,629	0	Gage Shale	MEMBER	P
MKD	2,712	0	Fort Riley Limestone	MEMBER	P
MKD	2,777	2,789	Florence Limestone	MEMBER	P
MKD	2,807	0	Wreford Limestone	FORMATION	P
MKD	2,832	0	Council Grove	GROUP	P
MKD	2,832	2,853.5	Speiser Shale	FORMATION	P
MKD	2,853.5	2,894.5	Funston Limestone	FORMATION	P
MKD	2,894.5	2,910.5	Blue Rapids Shale	FORMATION	P
MKD	2,910.5	2,929	Crouse Limestone	FORMATION	P
MKD	2,929	2,933.5	Easily Creek Shale	FORMATION	P
MKD	2,933.5	2,947	Middleburg Limestone	MEMBER	P
MKD	2,947	2,957.5	Hooser Shale	MEMBER	P
MKD	2,957.5	2,962.5	Eiss Limestone	MEMBER	P
MKD	2,962.5	2,972.5	Stearns Shale	FORMATION	P ▼
MKD	2,972.5	2,972.5	Merrill Limestone	MEMBER	P

Buttons: Clear Selection, Remove, Remove All, Load Data, Close

The MKD-07/2006 has only one top, so this dialog allows the user to add that top to the "User's Stratigraphic Units" Table. The MKD does not have this top and this is an extra top missing from the MKD data set.

KGS Stratigraphic Units:

☐ HUG ELOG-EM
☐ MKD
☒ MKD-07/2006

Add to User's Stratigraphic Units List:

☐ Remove & Replace ☒ Add to List ☐ Add New Units Only

Source	Top	Base	Name	Rank
MKD-07/2006	2,789	2,807	Matfield Shale	FORMAT

(1) Select the MKD-07/2006 Source, which will be displayed in the "Add to User's Stratigraphic Units List" Table.

(2) Select the "Add All" Button to move the contents from the "Add to User's Stratigraphic Units List" Table to the "User's Stratigraphic Units" table.

Add **Add All** Clear Selection

User's Stratigraphic Units:

Source	Top	Base	Name	Rank
MKD	2,538	2,580	Krider Limestone	MEMBER
MKD	2,629	0	Gage Shale	MEMBER
MKD	2,712	0	Fort Riley Limestone	MEMBER
MKD	2,777	2,789	Florence Limestone	MEMBER
MKD-07/2006	2,789	2,807	Matfield Shale	FORMATION
MKD	2,807	0	Wreford Limestone	FORMATION
MKD	2,832	0	Council Grove	GROUP
MKD	2,832	2,853.5	Speiser Shale	FORMATION
MKD	2,853.5	2,894.5	Funston Limestone	FORMATION
MKD	2,894.5	2,910.5	Blue Rapids Shale	FORMATION
MKD	2,910.5	2,929	Crouse Limestone	FORMATION
MKD	2,929	2,933.5	Easily Creek Shale	FORMATION
MKD	2,933.5	2,947	Middleburg Limestone	MEMBER
MKD	2,947	2,957.5	Hooser Shale	MEMBER
MKD	2,957.5	2,962.5	Eiss Limestone	MEMBER
MKD	2,962.5	2,972.5	Stearns Shale	FORMATION

Clear Selection

Remove

Remove All

Load Data

Close

Now with the data set complete select the "Load Data" Button to import the Tops data into the web app.

The Well data will automatically display to the "KIMELEON" dialog and load all the relevant panels, which can be accessed by the icon buttons at the top of the dialog. The "Horizon" icon button is the only panel that will show the Tops data with the summary of colorlith & color image plot tracks. Notice in the image below that the "K-U-Th" icon button is disabled, because there was no Spectral Gamma Ray Log Data present in the LAS File.

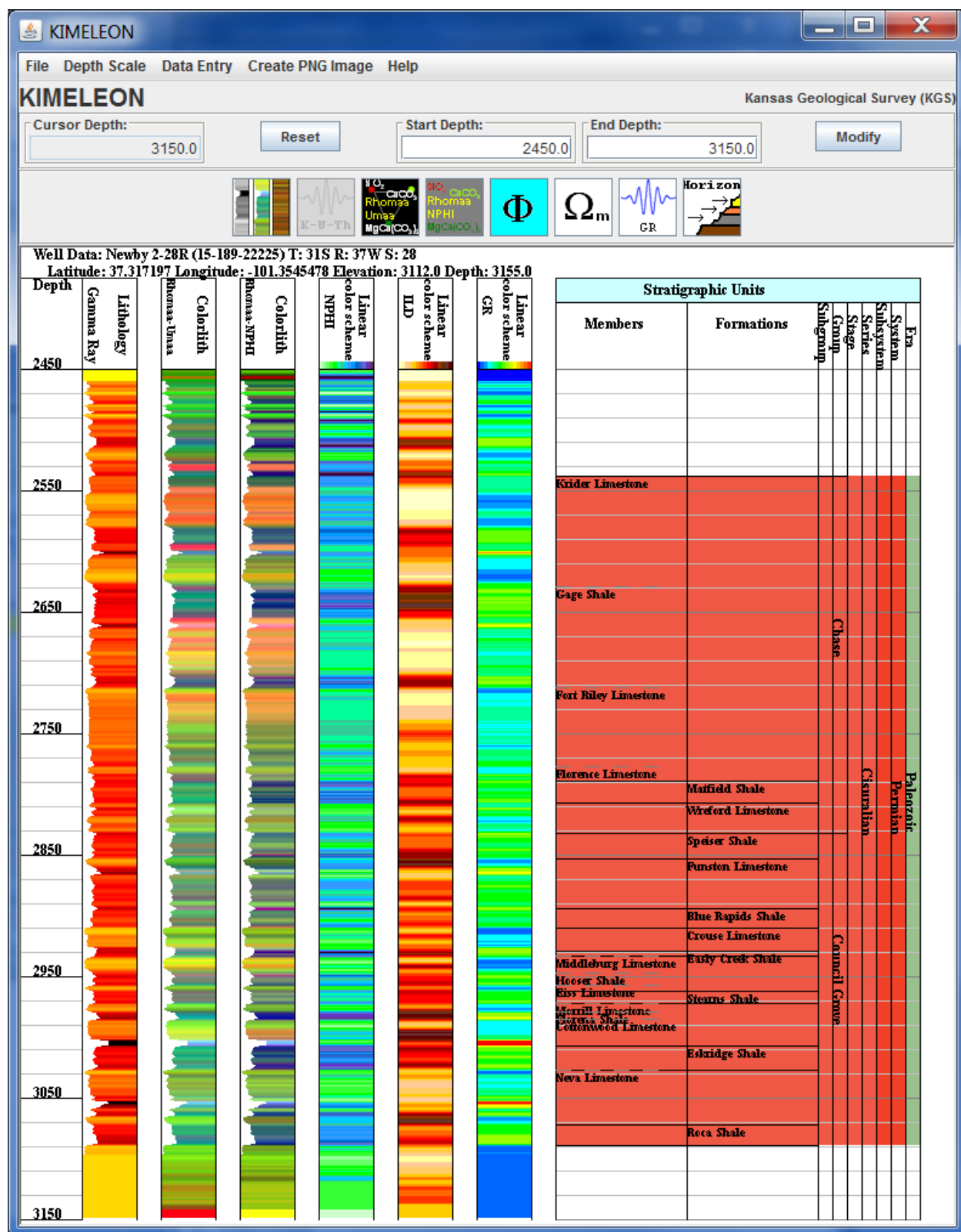


Figure: KIMELEON Dialog with the Newby 2-28R Well Data loaded from the KGS Database & Server.

Importing PC Data - Download Well Data to PC

Download either the ASCII Text Files directly or the Zip files extracting the contents into a directory. The problem with the ASCII Text Files being downloaded directly from a web page is that the web page will alter the contents so it does not retain the basic structure and add HTML text to the file. The preferred method if you have Zip or WinZip is to download the zip files to your PC and extract.

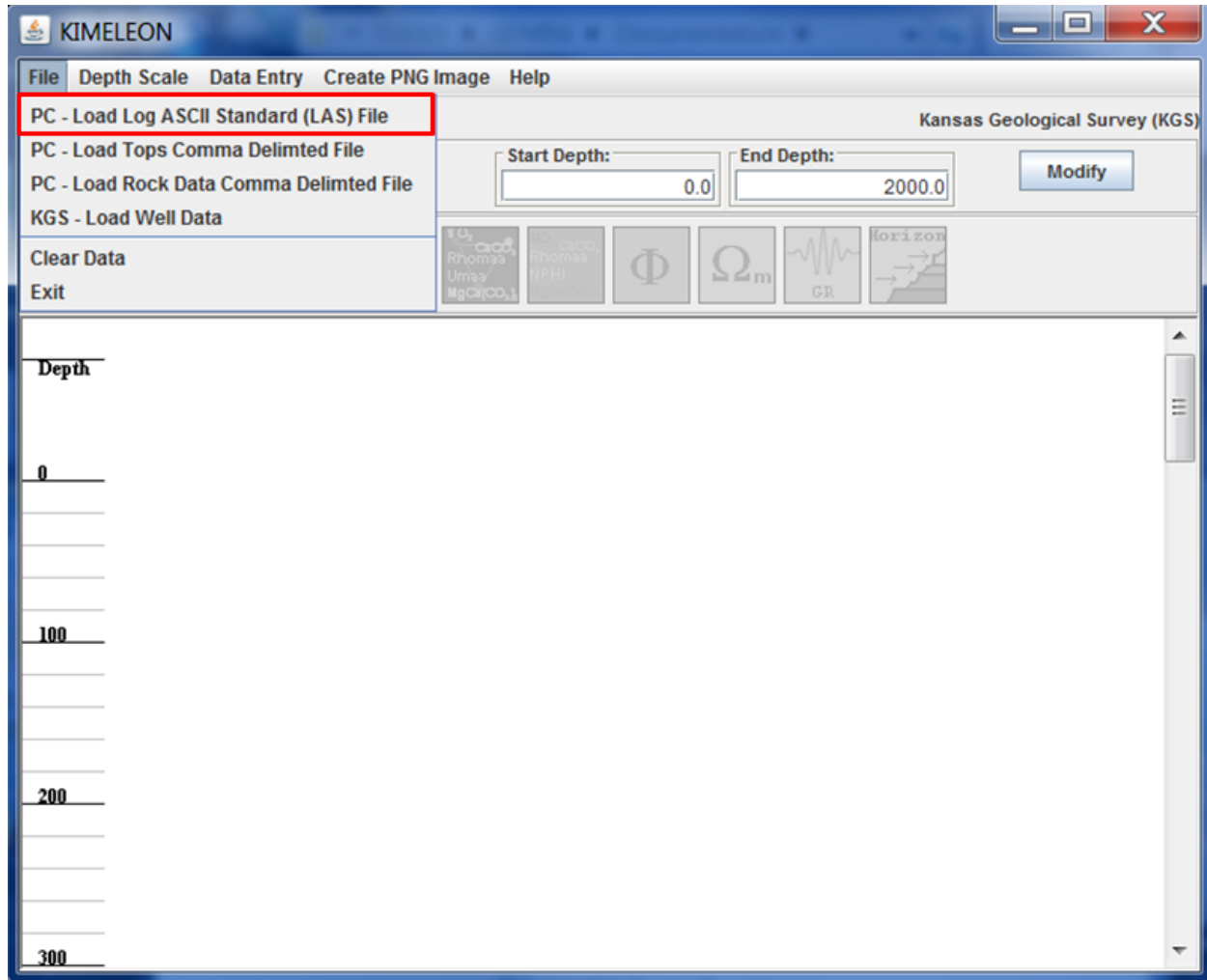
Well Data: Wellington KGS 1-32, Sumner County, Kansas

Type	ASCII Text Files
LAS 2.0	http://www.kgs.ku.edu/Gemini/Tools/documentation/Wellington-KGS-1-32.las
Tops	http://www.kgs.ku.edu/Gemini/Tools/documentation/Wellington-KGS-1-32_Tops.csv
Core	http://www.kgs.ku.edu/Gemini/Tools/documentation/Wellington-KGS-1-32_Core_Data.csv

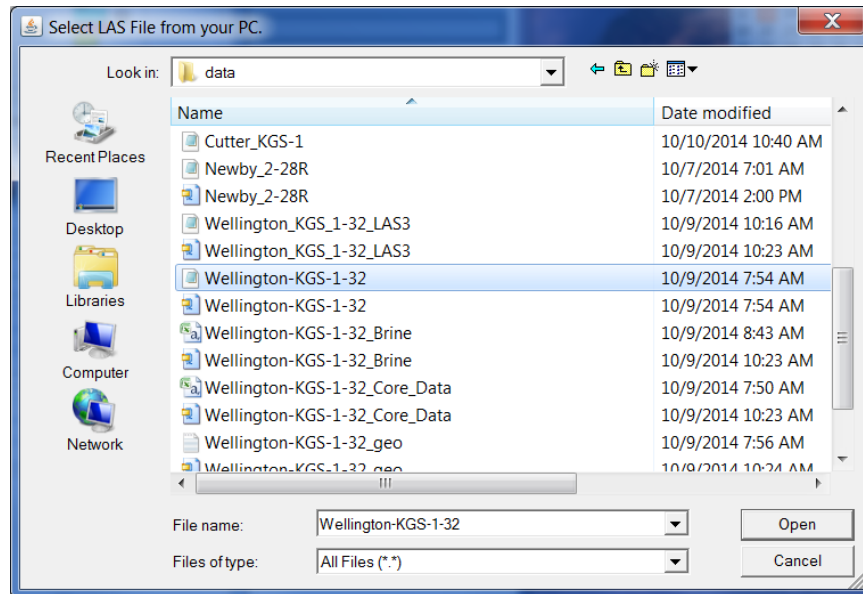
Type	Zip Files
LAS 2.0	http://www.kgs.ku.edu/Gemini/Tools/documentation/Wellington-KGS-1-32.zip
Tops	http://www.kgs.ku.edu/Gemini/Tools/documentation/Wellington-KGS-1-32_Tops.zip
Core	http://www.kgs.ku.edu/Gemini/Tools/documentation/Wellington-KGS-1-32_Core_Data.zip

Importing PC Data – Log ASCII Standard (LAS) version 2.0 File

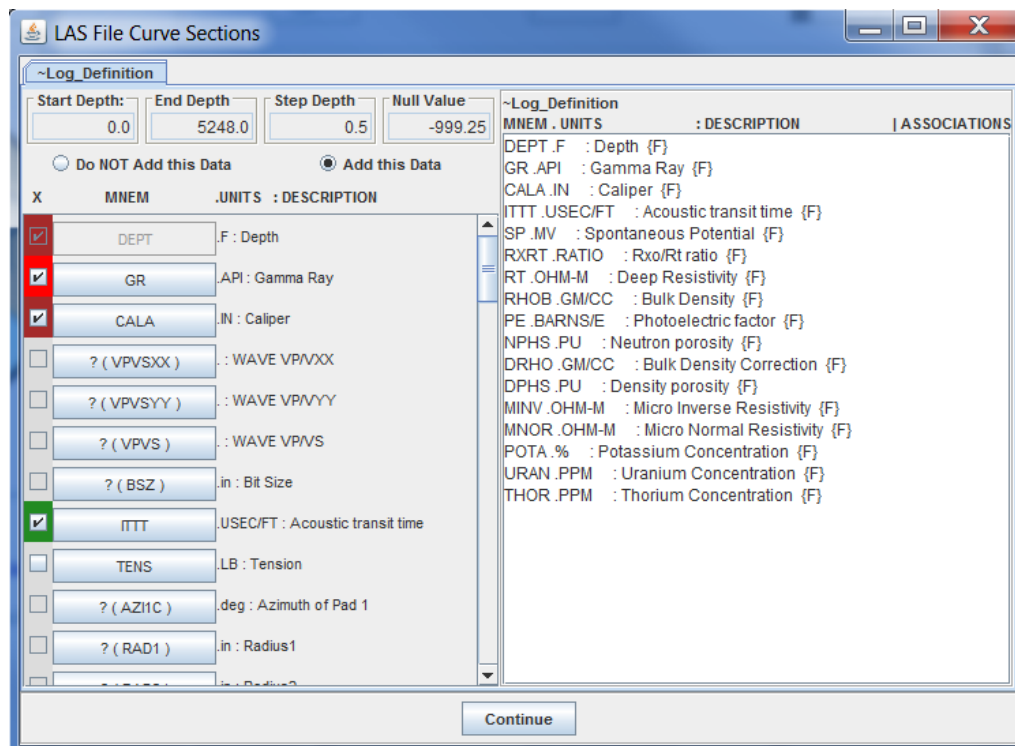
Most of the web apps will use the same input dialogs to import Log ASCII Standard (LAS) version 2.0 or 3.0 files. The Load Data Dialog is basically the same for most of the Web Apps, except they only load a subset of the total data types. In this example a LAS version 2.0 file is being imported into the web app.



Click on the “File” Menu in the menu bar at the top of the “KIMELEON” Dialog, which will display a list of menu options to import well data into the web app. Select the “PC - Load ASCII Standard (LAS) File” menu option. This will display the “Select LAS File from your PC” Dialog. This dialog allows the user to search their PC for the file of interest. In this example it is the LAS version 2.0 file Wellington-KGS-1-32.las, highlighted below. Select the Open button to display the “LAS File Curve Sections” Dialog.



The “LAS File Curve Sections” Dialog allows the user to map unknown LAS Curve Mnemonics to the KGS “Standard” Curve Mnemonics so they will be plotted in the LAS File Viewer Plot. This program reads the “LAS Tool Curve Mnemonics map to KGS Standard Mnemonics” XML File (http://www.kgs.ku.edu/software/gemini/data/las_standard_tools.xml), which will automatically maps the Curve Mnemonics from the LAS file to one of 31 KGS “Standard” Curve Mnemonics.



If a curve Mnemonic is not recognized the program will place a “?” in front of the Mnemonic, e.g. “?(BSZ)” for the “.in : Bit Size” Log Curve. If the user is satisfied with the automatic curve selections, which are checked and color coded, they only need to select the “Continue” Button at the bottom of the Dialog to import the file. The next section will take the user through a series of examples in changing the curve selections and mapping unknown curve mnemonics.

Notice that some of the check boxes are colored with different colors, which shows that the curves were automatically selected, but also to represent the curve type by color. The Curves are colored by type (data units) as follows,

- Orange - OHM-M or Resistivity Logs
- Cyan – PU or porosity Logs, Neutron Porosity, Density Porosity, etc.
- Greenish yellow – BARNS/E or Photoelectric Factor Logs
- Green – GM/CC or Bulk Density Log
- Forest Green – USEC/FT or the Acoustic Transit Time Log
- Red – API, PPM or “%” as Radioactive logs, Gamma Ray, Spectral Gamma Ray, etc.
- Blue – MD or Permeability Logs
- Brown – F, FT or IN or Depth
- Middle yellow – FRAC, or other log curve types.
- Dark Violet – UNI or Unknown Linear Curves
- Medium Violet – UNL or Unknown Logarithm Curves

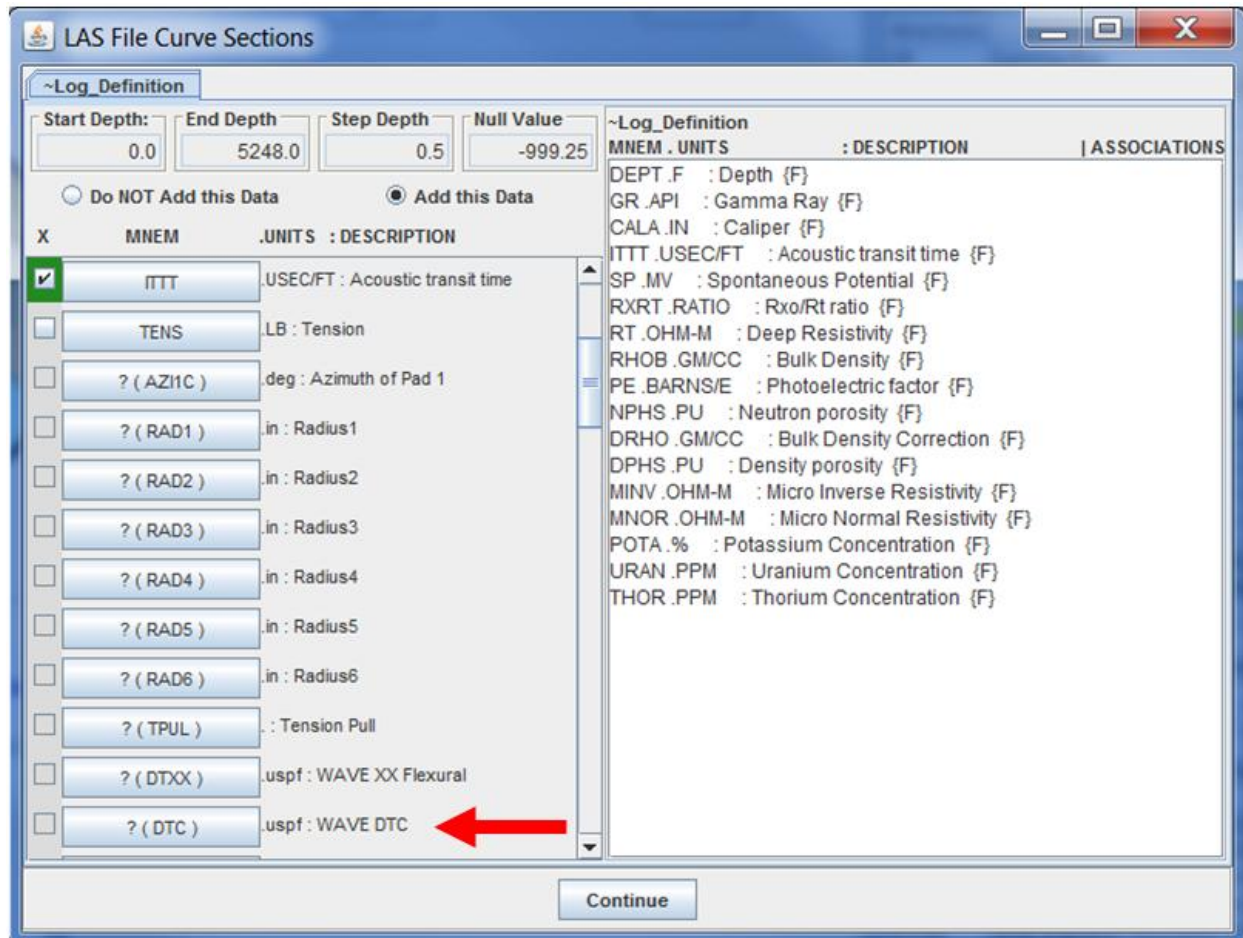
The color coding of the selected curves were added to also help the user visually recognize that a curve was selected or not.

Map Curves & Change Curve Selections

Some logs will have curve mnemonics that are not recognized as one of the KGS “Standard” Curve Mnemonics. The user will need to map the log curve to one of the KGS standard curves if they want to display the curve.

The first example is to map the Acoustic Transit Time (DT), which is labeled as “.uspf : WAVE DTC” log curve in the LAS File. Also notice that the button label “?(DTC)” is not recognized by the LAS FILE VIEWER web app.

Click on the “?(DTC)” Button to display the “Select KGS Standard Tools” Dialog. This dialog provides a list of the KGS “Standard” Curve Mnemonics, from which the user can map an unrecognized log curve to one of the KGS standard curve mnemonics. The KGS “Standard” Curve Mnemonics List was created as a way to standardize the alpha bit soup of Log Mnemonics.



Each logging company has their own curve mnemonics to represent similar tools. The LAS File Viewer program is a later version of code from the GEMINI Project LAS File Viewer Module, which needed to standardize the log curves so the curves could be automatically read and assigned a plot track. The “LAS Tool Curve Mnemonics map to KGS Standard Mnemonics” XML File was created to map the log curves from logs that were part of the KGS LAS File Collection which is not a complete list of possible curve mnemonics.

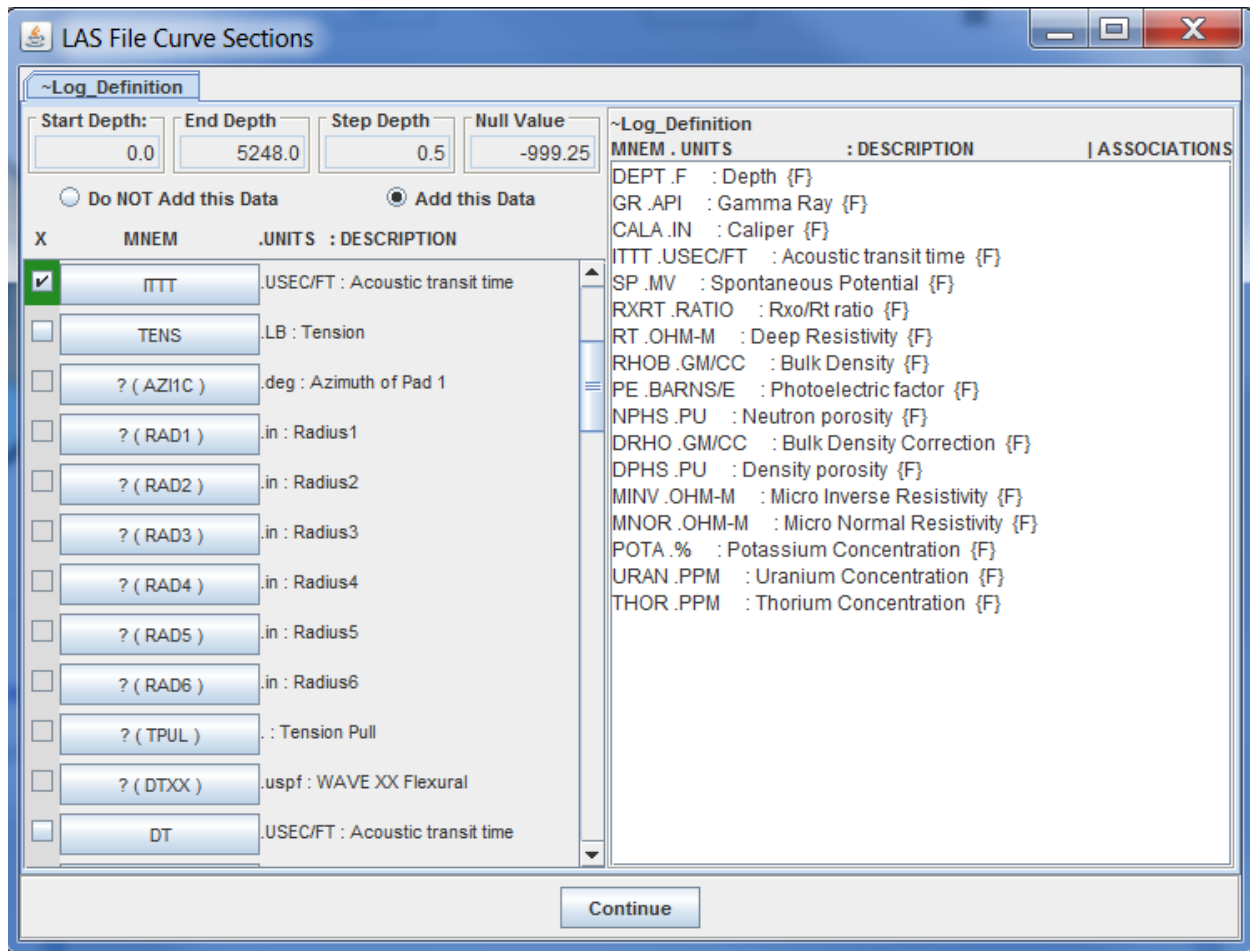
To map the unknown curve mnemonic “?(DTC)” you first notice that the unit is “uspf” (micro seconds per foot) a unit of time. Also the Acoustic Transit Time Curve Mnemonic is similar to the KGS “Standard” Curve Mnemonic “DT”. By selecting the “?(DTC)” Button you will display the “Select KGS Standard Tools” Dialog.

Select KGS Standard Tools

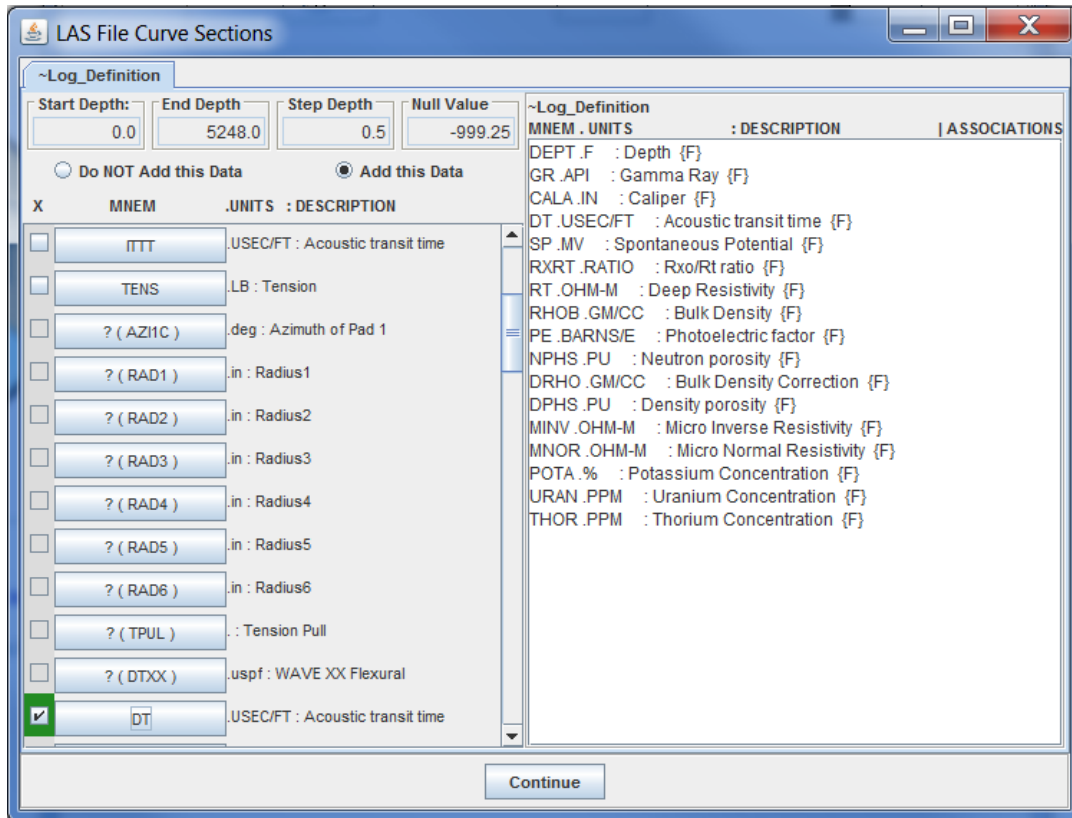
Mnemonic	Description	Units	Minimum	Maximum
GR	Gamma Ray	API	0	150
CGR	Gamma Ray Minus Uranium	API	0	150
SP	Spontaneous Potential	MV	0	1
CAL	Caliper	IN	6	12
PE	Photoelectric factor	BARN/E	0	20
RHOB	Bulk Density	GM/CC	2	3
DRHO	Bulk Density Correction	GM/CC	-1.5	0.5
DPHI	Density porosity	PU	-0.1	0.3
NPHI	Neutron porosity	PU	-0.1	0.3
SPHI	Sonic porosity	PU	-0.1	0.3
DT	Acoustic transit time	USEC/FT	40	140
COND	Conductivity	MMHO/M	0	2,000
CILD	Deep Induction Conductivity	MMHO/M	0	2,000
CILM	Medium Induction Conductivity	MMHO/M	0	2,000
RES	Resistivity	OHM-M	0.1	1,000
RDEP	Deep Resistivity	OHM-M	0.1	1,000
RMED	Medium Resistivity	OHM-M	0.1	1,000
RSHAL	Shallow Resistivity	OHM-M	0.1	1,000
ILD	Deep Induction Resistivity	OHM-M	0.1	1,000
ILM	Medium Induction Resistivity	OHM-M	0.1	1,000
SFLU	Spherically Focused Resistivity	OHM-M	0.1	1,000
LL	Deep Laterolog Resistivity	OHM-M	0.1	1,000
MLL	Micro Laterolog Resistivity	OHM-M	0.1	1,000
LL8	Shallow Laterolog Resistivity	OHM-M	0.1	1,000
LN	Long Normal Resistivity	OHM-M	0.1	1,000
SN	Shallow Normal Resistivity	OHM-M	0.1	1,000
MNOR	Micro Normal Resistivity	OHM-M	0.1	1,000
MSFL	Micro Spherically Focused Resistivity	OHM-M	0.1	1,000
MINV	Micro Inverse Resistivity	OHM-M	0.1	1,000
AHT10	Array Induction Resistivity-10	OHM-M	0.1	1,000
AHT20	Array Induction Resistivity-20	OHM-M	0.1	1,000
AHT30	Array Induction Resistivity-30	OHM-M	0.1	1,000
AHT60	Array Induction Resistivity-60	OHM-M	0.1	1,000
AHT90	Array Induction Resistivity-90	OHM-M	0.1	1,000
THOR	Thorium Concentration	PPM	-10	30

Select Cancel

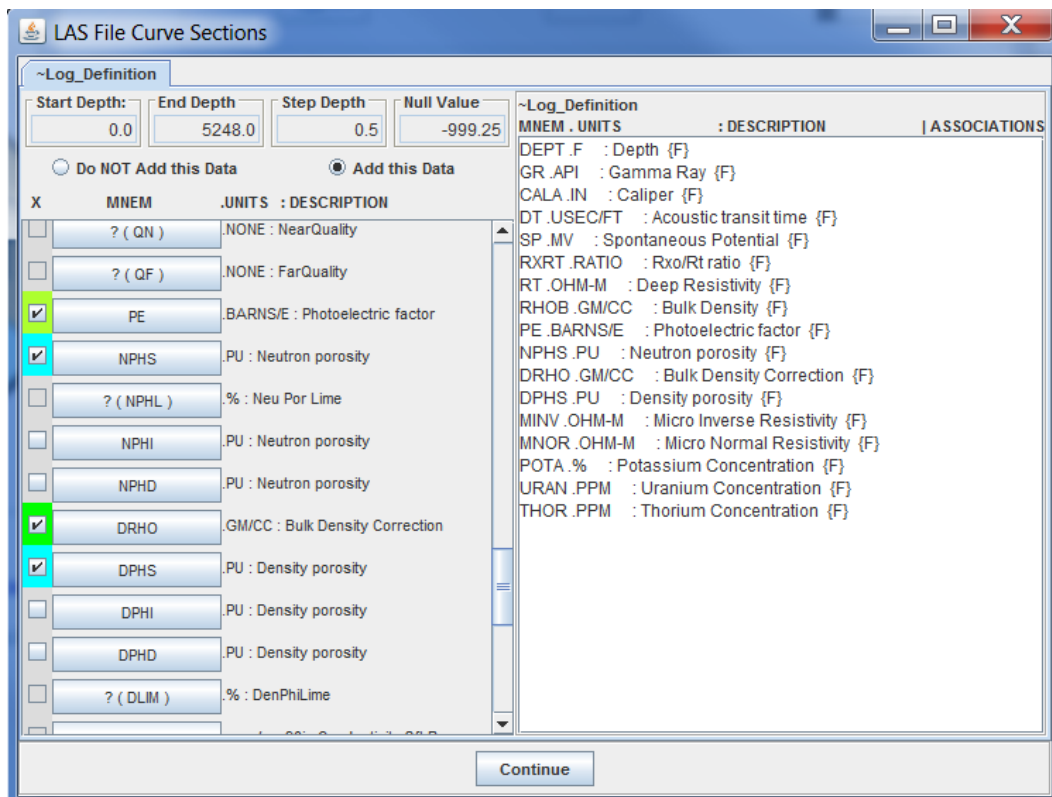
Highlight the “DT” Mnemonic Row and click on the “Select” Button to map the “?(DTC)” Curve Mnemonic to “DT” KGS Standard Curve Mnemonic.



The “?(DTC). .uspf : WAVE DTC” has changed to “DT.usc/ft : Acoustic transit time”. We want to change the selected “ITTT” Log Curve to “DT” Log Curve. The reason is that “ITTT” is the wrong curve type for the Acoustic Transit Time. The program found the curve mnemonic as similar to the “DT” Standard Curve Mnemonic, but this curve will not plot correctly in the LAS File Viewer Plot. Just click on the green check box in front of the “ITTT” Mnemonic Button to deselect the curve and then click on the check box in front of the “DT” Mnemonic Button to select it. Also notice that the ~Log_Definition Text Area was modified to show the change.



Moving the scroll bar down to the porosity curves, Neutron Porosity, and Density Porosity.



The LAS File Read will select the first curve that it recognizes and selects and color codes the curve. In this case the Neutron porosity mnemonic selected is “NPHS”, which is a valid curve, but the “NPHI” curve is desired so like the Acoustic Transit Time, you can deselect the “NPHS” and then select the “NPHI” Curve. Also the “DPHS” Density Porosity Curve can be deselected since the “RHOB” Bulk Density Curve has been selected. The reason for deselecting the Density Porosity Curve, if the Bulk Density Curve is present, is to force the LAS FILE VIEWER program to recompute the Density Porosity using a Limestone Matrix. If the Neutron Porosity, Bulk Density, Gamma Ray with/without a Photoelectric Factor Logs are present then the program will automatically compute a Lithology Composition Plot, but the Density Porosity has to be computed with a Limestone Matrix or the Lithology Composition Plot will not be computed correctly.

The screenshot shows the 'LAS File Curve Sections' dialog box. The 'Log Definition' tab is selected. At the top, there are input fields for 'Start Depth' (0.0), 'End Depth' (5248.0), 'Step Depth' (0.5), and 'Null Value' (-999.25). Below these are two radio buttons: 'Do NOT Add this Data' and 'Add this Data' (which is selected). The main area contains a table with columns 'X', 'MNEM', and '.UNITS : DESCRIPTION'. The table lists various curves, some of which are checked with a green box. The 'Continue' button is at the bottom right.

X	MNEM	.UNITS : DESCRIPTION
<input type="checkbox"/>	? (QN)	.NONE : NearQuality
<input type="checkbox"/>	? (QF)	.NONE : FarQuality
<input checked="" type="checkbox"/>	PE	.BARNS/E : Photoelectric factor
<input type="checkbox"/>	NPHS	.PU : Neutron porosity
<input type="checkbox"/>	? (NPHL)	.% : Neu Por Lime
<input checked="" type="checkbox"/>	NPHI	.PU : Neutron porosity
<input type="checkbox"/>	NPHD	.PU : Neutron porosity
<input checked="" type="checkbox"/>	DRHO	.GM/CC : Bulk Density Correction
<input type="checkbox"/>	DPHS	.PU : Density porosity
<input type="checkbox"/>	DPHI	.PU : Density porosity
<input type="checkbox"/>	DPHD	.PU : Density porosity
<input type="checkbox"/>	? (DLIM)	.% : DenPhilLime

The above dialog represents the changes made for the neutron/density porosity logs. The last curves to be modified are the Array Induction Logs. Haliburton uses a different curve mnemonic for these logs. Move the scroll bar up to find the Array Induction Logs, RT90, RT60, etc.

LAS File Curve Sections

~Log_Definition

Start Depth: 0.0 End Depth: 5248.0 Step Depth: 0.5 Null Value: -999.25

☐ Do NOT Add this Data ☒ Add this Data

X	MNEM	.UNITS : DESCRIPTION
<input checked="" type="checkbox"/>	SP	.MV : Spontaneous Potential
<input checked="" type="checkbox"/>	RXRT	.RATIO : Rxo/Rt ratio
<input type="checkbox"/>	RXO	.RATIO : Rxo/Rt ratio
<input type="checkbox"/>	? (RT90)	.ohmm : 90in Resistivity 2ft Res
<input type="checkbox"/>	? (RT60)	.ohmm : 60in Resistivity 2ft Res
<input type="checkbox"/>	? (RT30)	.ohmm : 30in Resistivity 2ft Res
<input type="checkbox"/>	? (RT20)	.ohmm : 20in Resistivity 2ft Res
<input type="checkbox"/>	? (RT10)	.ohmm : 10in Resistivity 2ft Res
<input checked="" type="checkbox"/>	RT	.OHM-M : Deep Resistivity
<input type="checkbox"/>	? (RMUD)	.ohmm : RMUD
<input checked="" type="checkbox"/>	RHOB	.GM/CC : Bulk Density
<input type="checkbox"/>	? (QN)	.NONE : NearQuality

~Log_Definition

MNEM	.UNITS	: DESCRIPTION	ASSOCIATIONS
DEPT.F		: Depth {F}	
GR.API		: Gamma Ray {F}	
CALA.IN		: Caliper {F}	
DT.USEC/FT		: Acoustic transit time {F}	
SP.MV		: Spontaneous Potential {F}	
RXRT.RATIO		: Rxo/Rt ratio {F}	
RT.OHM-M		: Deep Resistivity {F}	
RHOB.GM/CC		: Bulk Density {F}	
PE.BARNS/E		: Photoelectric factor {F}	
NPHI.PU		: Neutron porosity {F}	
DRHO.GM/CC		: Bulk Density Correction {F}	
MINV.OHM-M		: Micro Inverse Resistivity {F}	
MNOR.OHM-M		: Micro Normal Resistivity {F}	
POTA.%		: Potassium Concentration {F}	
URAN.PPM		: Uranium Concentration {F}	
THOR.PPM		: Thorium Concentration {F}	

Continue

Like the Acoustic Transit Time the “?(RT90).ohmm: 90in Resistivity 2ft Res” through “?(RT10).ohmm : 10in Resistivity 2ft Res” are not recognized. These curves can be map to the “AHT90 Array Induction Resistivity-90” to “AHT10 Array Induction Resistivity-10” KGS Curves Respectively. Click on the “?(RT90)” Mnemonic Button to display the “Select KGS Standard Tools” Dialog.

Select KGS Standard Tools

Mnemonic	Description	Units	Minimum	Maximum
GR	Gamma Ray	API	0	150
CGR	Gamma Ray Minus Uranium	API	0	150
SP	Spontaneous Potential	MV	0	1
CAL	Caliper	IN	6	12
PE	Photoelectric factor	BARNs/E	0	20
RHOB	Bulk Density	GM/CC	2	3
DRHO	Bulk Density Correction	GM/CC	-1.5	0.5
DPHI	Density porosity	PU	-0.1	0.3
NPHI	Neutron porosity	PU	-0.1	0.3
SPHI	Sonic porosity	PU	-0.1	0.3
DT	Acoustic transit time	USEC/FT	40	140
COND	Conductivity	MMHO/M	0	2,000
CILD	Deep Induction Conductivity	MMHO/M	0	2,000
CILM	Medium Induction Conductivity	MMHO/M	0	2,000
RES	Resistivity	OHM-M	0.1	1,000
RDEP	Deep Resistivity	OHM-M	0.1	1,000
RMED	Medium Resistivity	OHM-M	0.1	1,000
RSHAL	Shallow Resistivity	OHM-M	0.1	1,000
ILD	Deep Induction Resistivity	OHM-M	0.1	1,000
ILM	Medium Induction Resistivity	OHM-M	0.1	1,000
SFLU	Spherically Focused Resistivity	OHM-M	0.1	1,000
LL	Deep Laterolog Resistivity	OHM-M	0.1	1,000
MLL	Micro Laterolog Resistivity	OHM-M	0.1	1,000
LL8	Shallow Laterolog Resistivity	OHM-M	0.1	1,000
LN	Long Normal Resistivity	OHM-M	0.1	1,000
SN	Shallow Normal Resistivity	OHM-M	0.1	1,000
MNOR	Micro Normal Resistivity	OHM-M	0.1	1,000
MSFL	Micro Spherically Focused Resistivity	OHM-M	0.1	1,000
MINV	Micro Inverse Resistivity	OHM-M	0.1	1,000
AHT10	Array Induction Resistivity-10	OHM-M	0.1	1,000
AHT20	Array Induction Resistivity-20	OHM-M	0.1	1,000
AHT30	Array Induction Resistivity-30	OHM-M	0.1	1,000
AHT60	Array Induction Resistivity-60	OHM-M	0.1	1,000
AHT90	Array Induction Resistivity-90	OHM-M	0.1	1,000
THOR	Thorium Concentration	PPM	-10	30

Select Cancel

Highlight the AHT90 and click on the “Select” Button.

LAS File Curve Sections

~Log_Definition

Start Depth: 0.0 End Depth: 5248.0 Step Depth: 0.5 Null Value: -999.25

☐ Do NOT Add this Data ☒ Add this Data

X	MNEM	.UNITS : DESCRIPTION
<input checked="" type="checkbox"/>	SP	MV : Spontaneous Potential
<input checked="" type="checkbox"/>	RXRT	.RATIO : Rxo/Rt ratio
<input type="checkbox"/>	RXO	.RATIO : Rxo/Rt ratio
<input checked="" type="checkbox"/>	AHT90	.OHM-M : Array Induction Resistivity-90
<input type="checkbox"/>	? (RT60)	.ohmm : 60in Resistivity 2ft Res
<input type="checkbox"/>	? (RT30)	.ohmm : 30in Resistivity 2ft Res
<input type="checkbox"/>	? (RT20)	.ohmm : 20in Resistivity 2ft Res
<input type="checkbox"/>	? (RT10)	.ohmm : 10in Resistivity 2ft Res
<input checked="" type="checkbox"/>	RT	.OHM-M : Deep Resistivity
<input type="checkbox"/>	? (RMUD)	.ohmm : RMUD
<input checked="" type="checkbox"/>	RHOB	.GM/CC : Bulk Density
<input type="checkbox"/>	? (QN)	.NONE : NearQuality

~Log_Definition

MNEM . UNITS	: DESCRIPTION	ASSOCIATIONS
DEPT.F	: Depth {F}	
GR.API	: Gamma Ray {F}	
CALA.IN	: Caliper {F}	
DT.USEC/FT	: Acoustic transit time {F}	
SP.MV	: Spontaneous Potential {F}	
RXRT.RATIO	: Rxo/Rt ratio {F}	
AHT90.OHM-M	: Array Induction Resistivity-90 {F}	
RT.OHM-M	: Deep Resistivity {F}	
RHOB.GM/CC	: Bulk Density {F}	
PE.BARNS/E	: Photoelectric factor {F}	
NPHI.PU	: Neutron porosity {F}	
DRHO.GM/CC	: Bulk Density Correction {F}	
MINV.OHM-M	: Micro Inverse Resistivity {F}	
MNOR.OHM-M	: Micro Normal Resistivity {F}	
POTA.%	: Potassium Concentration {F}	
URAN.PPM	: Uranium Concentration {F}	
THOR.PPM	: Thorium Concentration {F}	

Continue

The “(?AHT90).ohmm : 90in Resistivity 2ft Res” has changed to “AHT90.OHM-M : Array Induction Resistivity-90” and the orange check box is selected. The rest of the Array Induction Log Curves each are mapped to the respective KGS Mnemonic Curve as follows,

(?RT90).ohmm : 90in Resistivity 2ft Res to AHT90.OHM-M : Array Induction Resistivity-90
 (?RT60).ohmm : 60in Resistivity 2ft Res to AHT60.OHM-M : Array Induction Resistivity-60
 (?RT30).ohmm : 30in Resistivity 2ft Res to AHT30.OHM-M : Array Induction Resistivity-30
 (?RT20).ohmm : 20in Resistivity 2ft Res to AHT20.OHM-M : Array Induction Resistivity-20
 (?RT10).ohmm : 10in Resistivity 2ft Res to AHT10.OHM-M : Array Induction Resistivity-10

LAS File Curve Sections

~Log_Definition

Start Depth: 0.0 End Depth: 5248.0 Step Depth: 0.5 Null Value: -999.25

☐ Do NOT Add this Data ☒ Add this Data

X	MNEM	.UNITS : DESCRIPTION
<input checked="" type="checkbox"/>	SP	.MV : Spontaneous Potential
<input checked="" type="checkbox"/>	RXRT	.RATIO : Rxo/Rt ratio
<input type="checkbox"/>	RXO	.RATIO : Rxo/Rt ratio
<input checked="" type="checkbox"/>	AHT90	.OHM-M : Array Induction Resistivity-90
<input checked="" type="checkbox"/>	AHT60	.OHM-M : Array Induction Resistivity-60
<input checked="" type="checkbox"/>	AHT30	.OHM-M : Array Induction Resistivity-30
<input checked="" type="checkbox"/>	AHT20	.OHM-M : Array Induction Resistivity-20
<input checked="" type="checkbox"/>	AHT10	.OHM-M : Array Induction Resistivity-10
<input checked="" type="checkbox"/>	RT	.OHM-M : Deep Resistivity
<input type="checkbox"/>	? (RMUD)	.ohmm : RMUD
<input checked="" type="checkbox"/>	RHOB	.GM/CC : Bulk Density
<input type="checkbox"/>	? (QN)	.NONE : NearQuality

~Log_Definition

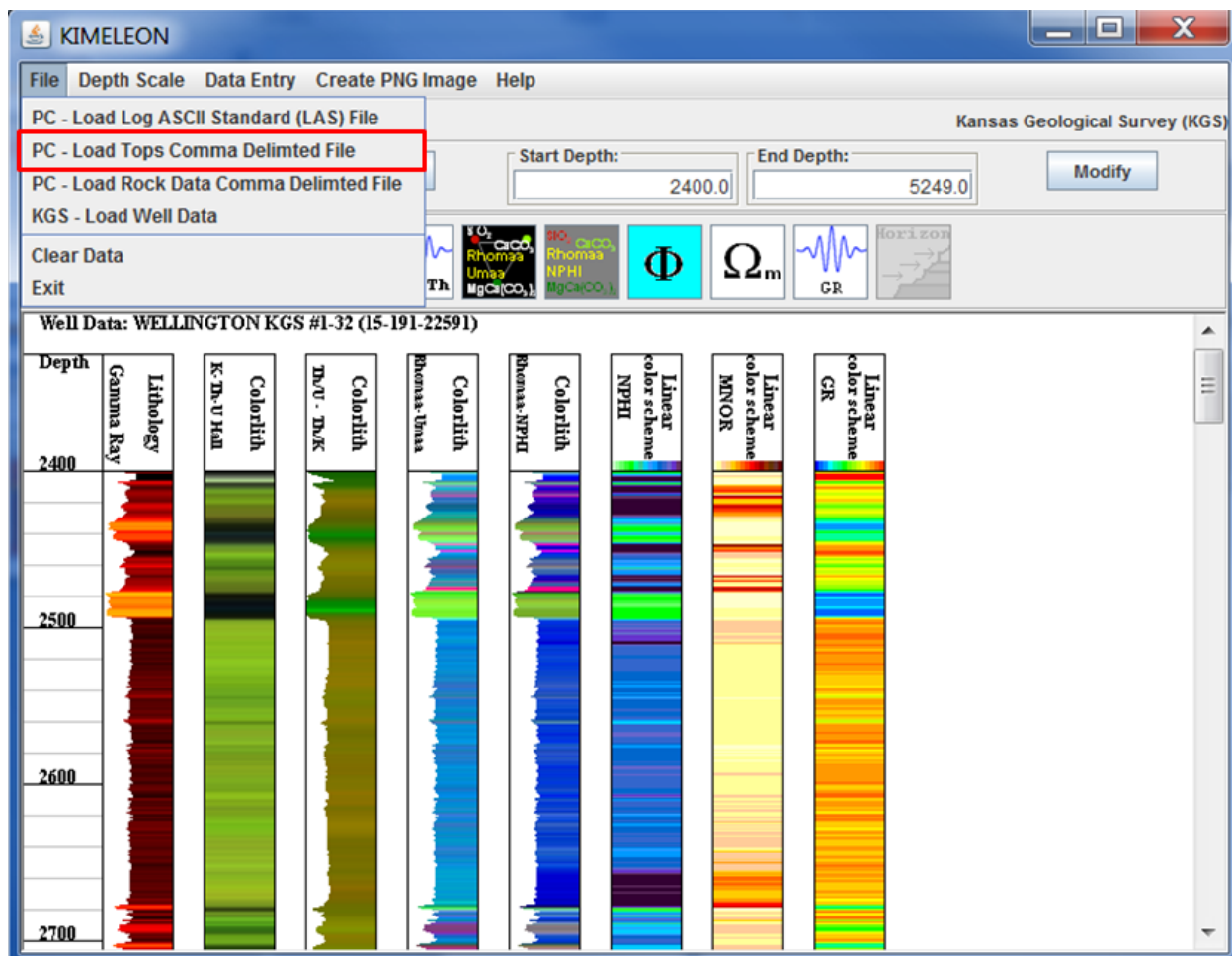
MNEM . UNITS	: DESCRIPTION	ASSOCIATIONS
DEPT .F	: Depth {F}	
GR .API	: Gamma Ray {F}	
CALA .IN	: Caliper {F}	
DT .USEC/FT	: Acoustic transit time {F}	
SP .MV	: Spontaneous Potential {F}	
RXRT .RATIO	: Rxo/Rt ratio {F}	
AHT90 .OHM-M	: Array Induction Resistivity-90 {F}	
AHT60 .OHM-M	: Array Induction Resistivity-60 {F}	
AHT30 .OHM-M	: Array Induction Resistivity-30 {F}	
AHT20 .OHM-M	: Array Induction Resistivity-20 {F}	
AHT10 .OHM-M	: Array Induction Resistivity-10 {F}	
RT .OHM-M	: Deep Resistivity {F}	
RHOB .GM/CC	: Bulk Density {F}	
PE .BARNSE	: Photoelectric factor {F}	
NPHI .PU	: Neutron porosity {F}	
DRHO .GM/CC	: Bulk Density Correction {F}	
MINV .OHM-M	: Micro Inverse Resistivity {F}	
MNOR .OHM-M	: Micro Normal Resistivity {F}	
POTA .%	: Potassium Concentration {F}	
URAN .PPM	: Uranium Concentration {F}	
THOR .PPM	: Thorium Concentration {F}	

Continue

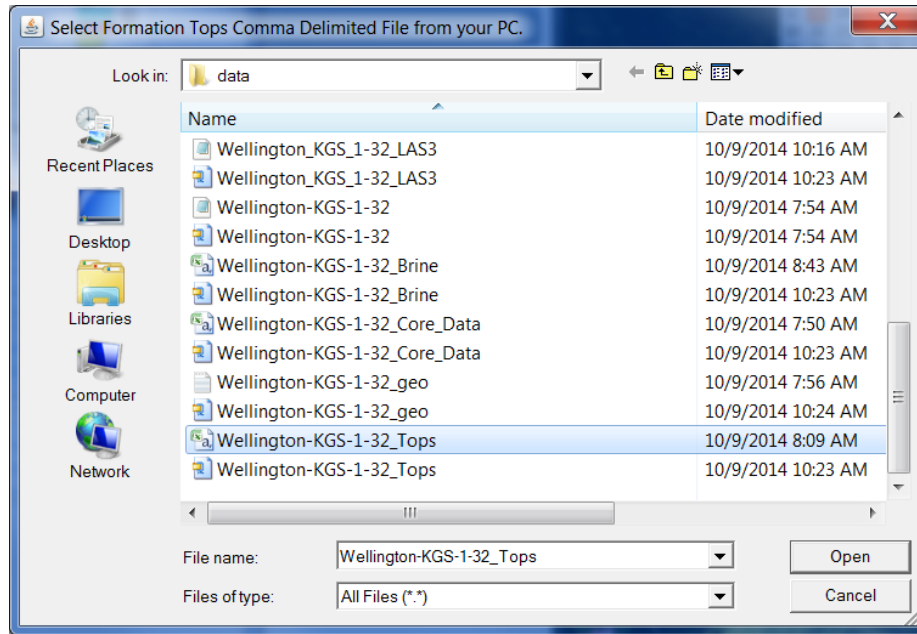
Select the Continue Button to read and parse the LAS log curves selected into the LAS FILE VIEWER Web App. Notice that the “Data Source Filenames:” Panel lists the LAS version 2.0 File that was just read in as well as the type of data, i.e. Log Data from LAS Data Type.

Importing PC Data – Tops CSV (Comma Separated Values) File.

Most of the web apps will use the same input dialogs to import tops CSV (Comma Separated Values) file. The Load Data Dialog is basically the same for most of the Web Apps except they only load a subset of the total data types. In this example a Tops CSV file is being imported into the web app.



Click on the “File” Menu in the menu bar at the top of the “KIMELEON” Dialog, which will display a list of menu options to import well data into the web app. Select the “PC - Load ASCII Standard (LAS) File” menu option. This will display the “Select Formation Tops Comma Delimited File from your PC” Dialog. This dialog allows the user to search their PC for the file of interest. In this example it is the Tops CSV file Wellington-KGS-1-32_Tops.csv, highlighted below. Select the Open button to display the “Map File Column Number to Region Column” Dialog.



The “Map File Column Number to Region Column” Dialog allows the user to map the file columns number to the web app tops data structure. In this example the file has the well information in line one of the Tops CSV File and line two of the Tops CSV File has the file data columns. In this case the chosen file columns match the Tops Mnemonics for the tops data structure. The File Column Number is automatically assigned to the Region Column Names. The user only needs to select the “Load Data” Button to parse the Tops Data into the web app.

Map File Column Number to Region Column

1st Line of Comma Delimited File:
Wellington KGS 1-32, 15-191-22591, T31S R1W sec. 32, GL:1259, KB:1272

2nd Line of Comma Delimited File:
Top, Name, Rank, System, Subsystem, Series, source

Formation Tops Columns:
Start Reading Data at Row Assume Row & Column Count is 1,2,3 ...

Region Column Name	File Column Number
Depth Top	<input type="text" value="1"/>
Depth Base	<input type="text" value="0"/>
Stratigraphic Unit Rank [SYSTEM, GROUP, etc.]	<input type="text" value="3"/>
Stratigraphic Name	<input type="text" value="2"/>
Alternate Name	<input type="text" value="0"/>
Era	<input type="text" value="0"/>
System	<input type="text" value="4"/>
Subsystem	<input type="text" value="5"/>
Series	<input type="text" value="6"/>
Subseries { Pennsylvanian & Mississippian Series }	<input type="text" value="0"/>
Stage	<input type="text" value="0"/>
Group	<input type="text" value="0"/>
Subgroup	<input type="text" value="0"/>
Formation	<input type="text" value="0"/>
Start Age (Ma)	<input type="text" value="0"/>
End Age (Ma)	<input type="text" value="0"/>

Tops CSV (Comma Separated Values) File Structure.

The Wellington KGS 1-32 Tops CSV example has two introduction lines, the first line is the well header information and the second line is the actual column labels for the tops data, illustrated below,


```

Line 1 Well Header Info Wellington KGS 1-32, 15-191-22591, T31S R1W sec. 32, GL:1259, KB:1272
Line 2 Data Column Labels Top, Name, Rank, System, Subsystem, Series, source
Line 3 Data Start 620, Chase, GROUP, Permian, , Wolfcampian, PG
748, Towanda Limestone, MEMBER, Permian, , Wolfcampian, PG
1595, Wabaunsee, GROUP, Carboniferous, Pennsylvanian, Upper, PG
1622, Root Shale, FORMATION, Carboniferous, Pennsylvanian, Upper, PG
1662, Stotler Limestone, FORMATION, Carboniferous, Pennsylvanian, Upper, PG
1920, Severy Shale, FORMATION, Carboniferous, Pennsylvanian, Upper, PG
1980, Topeka Limestone, FORMATION, Carboniferous, Pennsylvanian, Upper, PG
2312, Lecompton Limestone, FORMATION, Carboniferous, Pennsylvanian, Upper, PG
2402, Heebner Shale, MEMBER, Carboniferous, Pennsylvanian, Upper, PG
2703, Stalnaker Sandstone, BED, Carboniferous, Pennsylvanian, Upper, PG
3039, Kansas City, GROUP, Carboniferous, Pennsylvanian, Upper, PG
3169, Stark Shale, MEMBER, Carboniferous, Pennsylvanian, Upper, PG

```

Figure: Partial Contents of the Wellington-KGS-1-32_Tops.csv File.

The “Map File Column Number to Region Column” Dialog allows the user to map the data in the Tops CSV File to the web app data structure variables. The program first reads the first and second line of the CSV File looking for the data column headers. The lines are each parsed to single out the data column headers and to match those headers to the tops data structure. The program then assigns the column number to the Region Column Name starting at column 1,2,3, ... if the file column name used matches the expected region column name. The Column Names matrix used to parse the file column variables are listed below,

Depth Top	Top	Start
Depth Base	Base	End
Stratigraphic Unit Rank [SYSTEM, GROUP, etc.]	Rank	
Stratigraphic Name	Name	
Alternate Name	Alt Name	
Era		
System	Sys	
Subsystem	subsys	
Series	Ser	
Subseries { Pennsylvanian & Mississippian Series }	Subseries	Subser
Stage	Stg	
Group	Grp	
Subgroup	subgrp	
Formation	Form	
Start Age (Ma)	Start Age	
End Age (Ma)	End Age	

The Wellington KGS 1-32 Tops CSV File example above line 2 has only the Top, Tops Name, Rank, System, Subsystem, Series and Source as the column name variables. The program was able to map each of the column headers to the tops data structure, except Source, i.e.

Column	File Column Label	Tops Data Name
1	Top	Depth Top
2	Name	Stratigraphic Name
3	Rank	Stratigraphic Unit Rank
4	System	System
5	Subsystem	Subsystem
6	Series	Series
7	Source	

KIMELEON

File Depth Scale Data Entry Create PNG Image Help

Kansas Geological Survey (KGS)

Cursor Depth: 3116.0 Reset Start Depth: 2400.0 End Depth: 5249.0 Modify

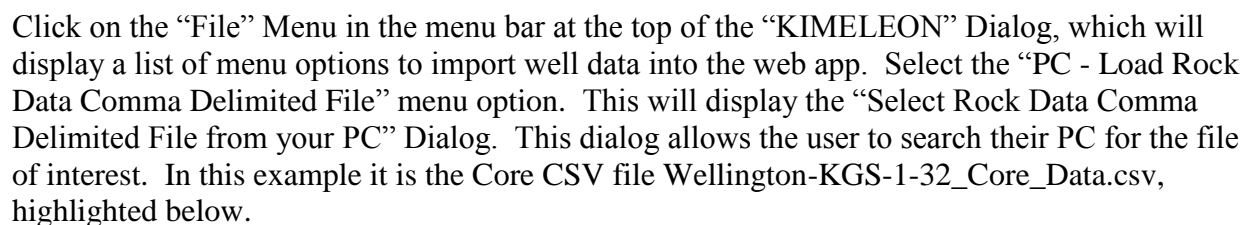
Well Data: WELLINGTON KGS #1-32 (15-191-22591)

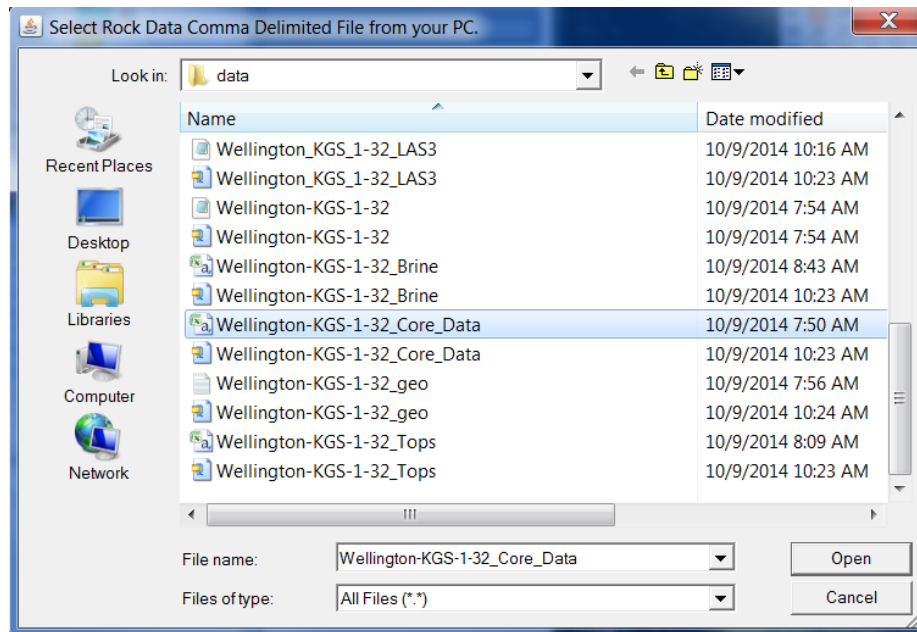
Depth	Lithology Gamma Ray	Colorith K-Th-U-Mn	Colorith Th/U - D/K	Colorith Rhomax Umas	Colorith Rhomax NPHI	Linear color scheme NPHI	Linear color scheme MINOR	Linear color scheme GR
2400								
2500								
2600								
2700								
2800								
2900								
3000								
3100								

Stratigraphic Units						
Members	Formations	Subgroup	Stage	Series	System	Era
Hedner Shale						
Stalnaker Sandstone						
				Upper	Pennsylvanian	

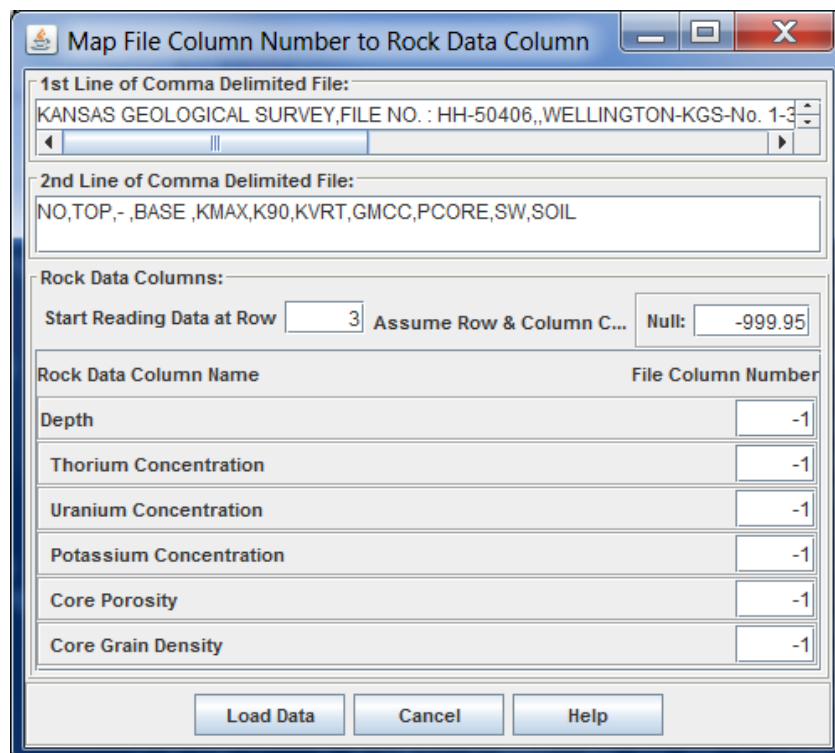
37

Most of the web apps will use the same input dialogs to import Core Data CSV (Comma Separated Values) file. The Load Data Dialog is basically the same for most of the Web Apps, except they only load a subset of the total data types. In this example a Core Data CSV file is being imported into the web app.





Select the Open button to display the “Map File Column Number to Rock Column” Dialog.



The “Map File Column Number to Rock Data Column” Dialog allows the user to map the file columns number to the web app core data structure. This dialog maps imports the core data and

loads the data into the LAS File Data Structure so the curves can be plotted as color image and/or colorlith plot tracks.

Core Data CSV (Comma Separated Values) File Structure.

The Wellington KGS 1-32 Core Data CSV example has two introduction lines, the first line is the well header information and the second line is the actual column labels for the core data, illustrated below,

```

Line 1 Well Header Info KANSAS GEOLOGICAL SURVEY,FILE NO. : HH-50406,,WELLINGTON-KGS-No. 1-32,FIELD : WELLINGTON,
Line 2 Data Column Labels NO,TOP,- ,BASE ,KMAX,K90,KVRT,GMCC,PCORE,SW,SOIL
Line 3 Data Start 2-29,3627.55,-,3628.00,0.02,<.01,<.01,2.69,0.3,90.9,0.0
2-39,3638.40,-,3638.90,0.01,<.01,<.01,2.70,0.6,80.5,0.0
3-0,3660.40,-,3660.75,2.00,0.08,2.29,2.62,4.7,82.5,0.0
3-1,3661.70,-,3662.00,2.72,1.74,0.02,2.61,6.8,75.1,0.0
3-2,3662.50,-,3663.00,3.70,2.96,4.55,2.59,11.1,45.5,22.5
3-3,3663.00,-,3663.50,25.36,9.94,29.36,2.60,14.1,41.2,29.2
3-4,3664.30,-,3664.75,12.42,8.97,3.84,2.62,8.0,97.6,0.0
3-5,3665.00,-,3665.60,4.20,3.40,3.38,2.61,7.4,98.6,0.0
3-6,3666.00,-,3666.25,8.64,0.40,<.01,2.58,5.5,46.0,10.7
3-7,3667.20,-,3667.70,6.56,5.97,1.93,2.60,5.8,93.9,0.0
3-8,3668.50,-,3669.00,21.77,10.68,5.78,2.59,12.0,82.4,0.0
3-9,3669.30,-,3669.80,20.30,20.03,10.96,2.60,11.8,70.4,0.0
3-10,3670.00,-,3670.60,64.64,60.27,35.99,2.61,13.1,75.8,0.0
3-11,3671.00,-,3671.50,9.68,8.11,2.41,2.79,19.7,50.2,23.9
3-12,3672.30,-,3672.80,14.08,13.60,7.96,2.77,22.7,48.0,27.1
3-13,3673.40,-,3673.90,13.37,12.95,8.20,2.79,24.3,50.1,24.4
3-14,3674.00,-,3674.50,21.85,21.24,9.30,2.77,22.9,52.8,28.2
3-15,3675.30,-,3675.80,22.31,19.92,11.61,2.78,26.4,58.8,25.2

```

Figure: Partial Contents of the Wellington-KGS-1-32_Core_Data.csv File.

The “Map File Column Number to Rock Data Column” Dialog allows the user to map the data in the Core Data CSV File to the web app data structure variables. The program first reads the first and second line of the CSV File looking for the data column headers. The lines are each parsed to single out the data column headers and to match those headers to the core data structure. The program then assigns the column number to the Rock Data Column Name starting at column 1,2,3, ... if the file column name used matches the expected rock data column name. The Column Names matrix used to parse the file column variables are listed below,

Mnem	Description
Depth	Depth Top
THOR	Thorium Concentration
URAN	Uranium Concentration
POTA	Potassium Concentration
PHI	Core Porosity
GMCC	Grain Density

The Wellington KGS 1-32 Core Data CSV File example above line 2 has only the Top, Base KMAX, K90, KVRT, GMCC, PCORE, SW and SOIL as the column name variables. The user will only be able to map a limited number of column headers to the core data structure, i.e.

Column	File Column Label	Core Data Name
1	NO	
2	TOP	Depth
3	-	
4	BASE	
5	KMAX	
6	K90	
7	KVRT	
8	GMCC	Grain Density (gm/cc)
9	PCORE	Whole Core Porosity
10	SW	
11	SOIL	

Map File Column Number to Rock Data Column

1st Line of Comma Delimited File:
KANSAS GEOLOGICAL SURVEY,FILE NO. : HH-50406,,WELLINGTON-KGS-No. 1-3

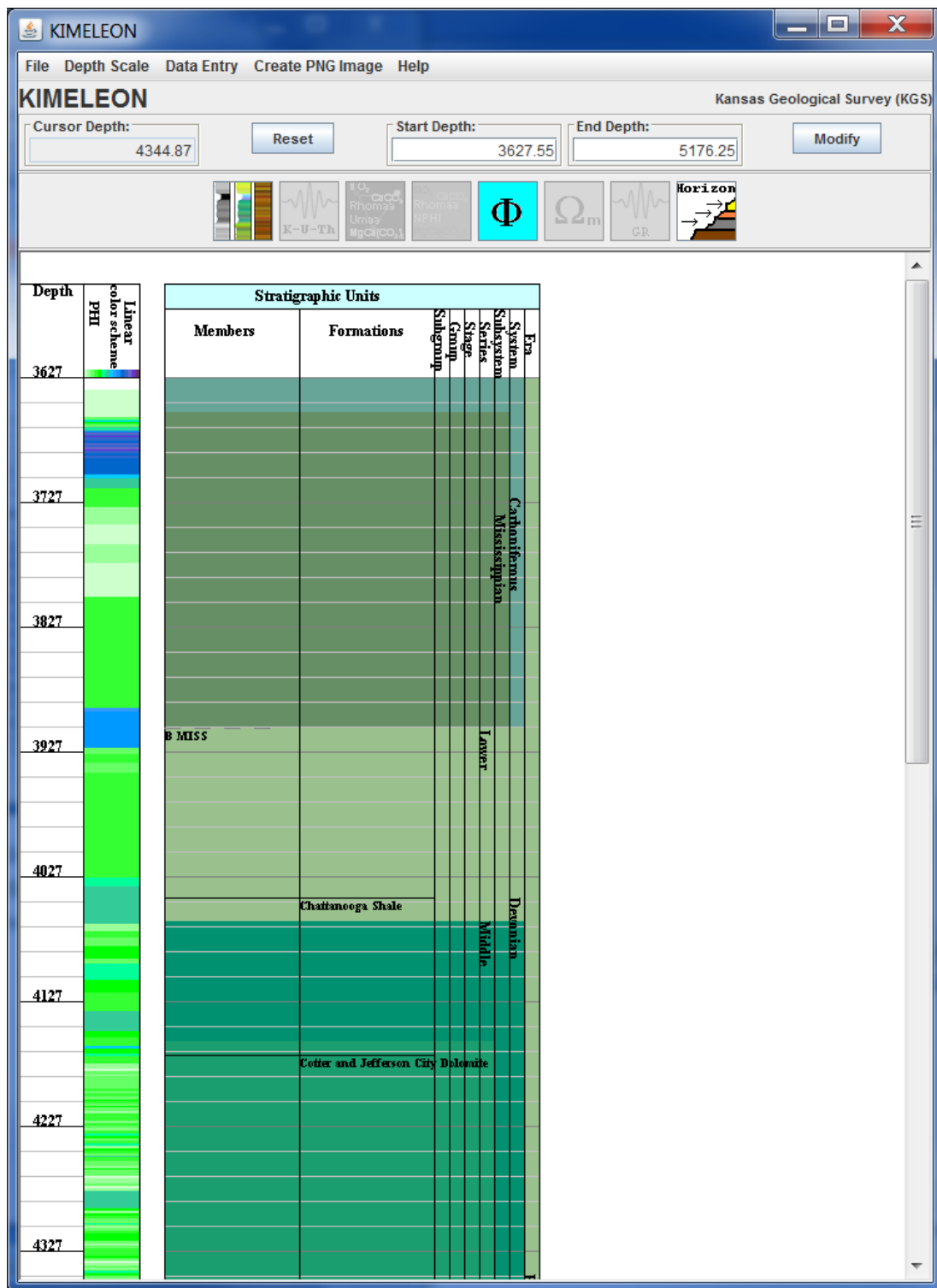
2nd Line of Comma Delimited File:
NO,TOP,-,BASE,KMAX,K90,KVRT,GMCC,PCORE,SW,SOIL

Rock Data Columns:

Start Reading Data at Row Assume Row & Column C... Null:

Rock Data Column Name	File Column Number
Depth	<input type="text" value="2"/>
Thorium Concentration	<input type="text" value="-1"/>
Uranium Concentration	<input type="text" value="-1"/>
Potassium Concentration	<input type="text" value="-1"/>
Core Porosity	<input type="text" value="9"/>
Core Grain Density	<input type="text" value="8"/>

When the user selects the “Load Data” Button on the “Map File Column Number to Rock Data Column” Dialog the data is parsed into the LAS File Viewer Program, where the Core Data CSV file name is entered into the “PC ASCII Files:” Panel as well as the data type source. The user can import the tops CSV file as a reference to the porosity core data.



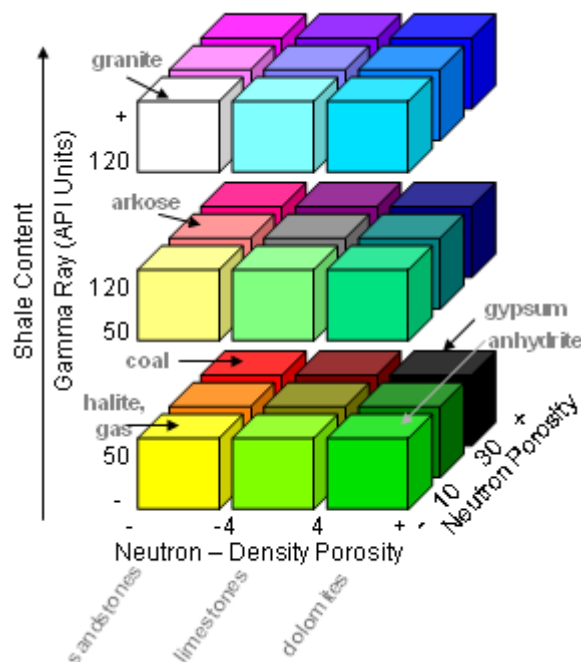
The Well data will automatically display to the “KIMELEON” dialog and load all the relevant panels, which can be accessed by the icon buttons at the top of the dialog. The “Horizon” icon button is the only panel that will show the Tops data with the summary of colorlith & color image plot tracks.

KIMELEON History

Peter L. Briggs⁽¹⁾ identified a method to assist geologist in well log interpretation by creating a color log presentation. His method provided a means to combine three different log curves from one well into one image track that varies along the depth by assigning each curve to a specific primary color of red, green and blue. His method assumed that the primary colors would appear to the human eye as orthogonal and miscible where the resulting color image would preserve all the information of the original log curves. The color image presentation of well log data reduced the number of displays competing for the interpreter's attentions by uniting the three log curves into one display and relieved the burden of mentally combining data from several separate displays. Overall the color log images presented log data to the user in a form that differs from the conventional log presentation and relied on the human perceptions and pattern recognition skills.



David R. Collins & John H. Doveton⁽²⁾ extended this technique to identify lithology by using the neutron & density porosity and gamma ray logs, by noting that several porosity log readings gave direct indications of the rock mineralogy. The neutron and density logs were first overlaid on a common scale of equivalent limestone porosity units and their relative disposition considered together with the gamma ray log. Shales have high neutron porosity (due to bound water), relatively low density reading and a high gamma ray value. Limestones have low gamma ray value, and a coincident neutron and density porosity. Dolomites have a low gamma ray value, a relatively low density reading and a relatively high neutron reading. Sandstones have a low gamma ray value, a relatively high density reading and relatively low neutron reading.



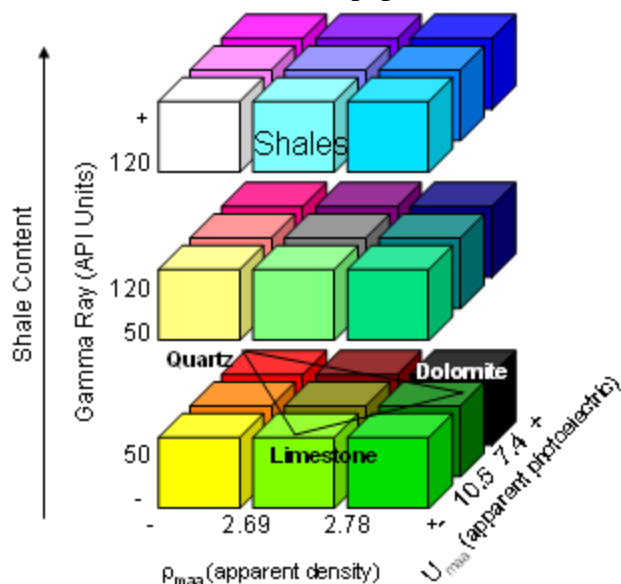
They based the lithology identification on three decision rules:

1. Is the gamma ray reading low, intermediate or high?
2. Is the density reading higher than, approximately equal to, or less than the neutron reading?
3. Do the density and neutron readings indicate a low, intermediate, or high porosity?

They extended the 27 possible responses to a color cube whose orthogonal axes match the rules. The design rules of the cube were geared to a basic discrimination between the potential reservoir lithology of sandstone, limestone, and dolomite, and their distinction from shale. They also noted that other lithologies have characteristic log responses which made additional identifications possible.



David Collins⁽³⁾ in a USGS paper discusses the visualization of subsurface geology as achieved

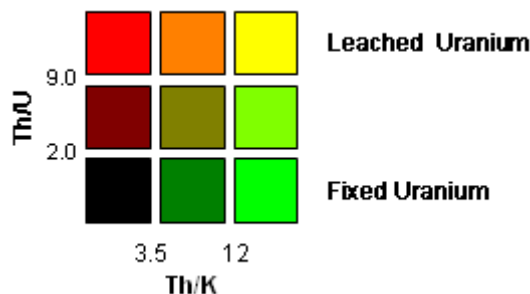


by COLORLITH, a software system developed at the Kansas Geological Survey in an effort to provide low-cost, high-resolution, interpretation, and visualization of well log data from a single well or multiple wells. The COLORLITH system was designed to run on UNIX workstation environments written in FORTRAN77. The input was digital LAS log data and the output was written to a PostScript file. COLORLITH was specifically designed to analyze and display gamma ray and litho-density logs. The COLORLITH computes the apparent density (Rhomaa) and the apparent photoelectric factor (Umaa) from the litho-density logs, Photoelectric factor, Bulk Density and the Neutron Porosity log curves.

KIMELEON extends both the Porosity Difference Method and the Rhomaa-Umaa Method that is computed in COLORLITH system. Both of the earlier versions done by Doveton & Collins^(2,3) use a 3 X 3 X 3 color cube to create a color image track. KIMELEON extends the color cube to 256 X 256 X 256 colors to create the color image track which increases the sensitivity to the log responses and will bring out more subtle changes in the lithology over the depth. This new system is web based and is written in JAVA and allows the user to directly interact with the data in real time. The digital LAS files can be retrieved from the user's PC or from the KGS server and plotted to a well profile plot.



Colorlith has been developed over the past 20 years from Briggs to Doveton & Collins and has been defined by Schlumberger as "A system for color-coding three-dimensional information. This system is used in wire line log analysis to provide color shading in which the final color is determined by the values of three curves. One curve dictates the intensity of red, a second the intensity of green, and the third the intensity of blue. The final resulting color is the result of the three input curves. The input curves may be raw curves from the field or computed curves. When used for correlation work on cross sections, the curves must have been normalized to remove the effects of incorrect calibrations and borehole problems." The colorlith or Briggs Cube is considered an excellent method for condensing the wire line log curves into one track to visualize the lithologies of a well. Another method used and developed by Doveton⁽⁴⁾ is the spectral gamma ray or Th/U and Th/K ratio cross plots introduced with the KGS #1 Braun in Ellis County Kansas.



The spectral gamma ray log can be used to estimate volumes and types of clay minerals, and is also useful for identifying fractures that have uranium salts precipitated in them by ground-water. Significant potassium and thorium concentrations in carbonates are also found in clay minerals and may show on porosity logs as shales. The spectral gamma-ray log helps to differentiate the radioactive carbonates from shales and clays. Several case studies of the use of spectral gamma-ray logs in Kansas show how the data can be cross-plotted using digital recording applications to identify facies. Natural gamma radiation in rocks is almost entirely attributable to potassium-40 and the radioactive isotopes of the uranium and thorium families. The concentrations of the three main radioactive elements Potassium, Uranium and Thorium in the formation can often be used to give an indication of the mineralogy and/or geochemistry. Thorium may be associated with an increase of terrigenous clays. Uranium is frequently associated with the presence of organic matter in black shale deposits. In sandstone, high Potassium may be caused by the presence of potassium feldspars or micas.

Th/K	Mineral
0.5 to 0.6	Feldspars
0.6 to 2.0	Micas
2.0 to 3.5	Illite
3.5 to 12.0	Smectite
12.0 to 28.0	Kaolinite-Chlorite
28.0+	Heavy Th-Bearing Minerals



Another colorlith technique proposed by Matt Hall⁵ uses the Spectral Gamma- Ray logs, Potassium (red), Thorium (green) and Uranium (blue). His method maps the gamma ray spectrum to the visible end of the electromagnetic spectrum. The gamma-ray emissions from the decay of potassium-40 nuclei have the lowest energy, the potassium log is represented by shades of red, thorium by green, and uranium by blue. He noted that logs are represented by monochrome variable density displays. Low relative amplitudes translate to low color values and are displayed as dark colors, and high amplitudes are displayed as bright, saturated colors. The colors are then combined additively in a RGB spectrum producing a composite image.

The "micro-electrical imaging" technique replaces the original current measurement of the electrodes from a high precision form to normalized, quantization form with the measurements being significantly reduced. The purpose is to produce an image with color levels and contrasts that make it easier for the interpreter to analyze. Colors are represented as dark have been chosen to represent high conductivities because shales are often more conductive and darker-colored than other lithologies.



Normalization of the image over a given depth range can enhance the contrast of the image. The number of data points over the given depth range is subdivided equally into each color bin. This is the non-linear process and a great degree contrast that helps spotting image details in areas where sample values differ little from each other. These images often appear somewhat harsh in contrast.



Quantization procedure is very similar to dynamic normalization, but generally the assignment of colors is made using the dynamics of the entire length of the log. The number of data points is distributed by magnitude into equal bins. This often results in an image where intermediate colors abound and few extreme colors such as dark brown or cream occur, except for unusual distribution of the input data. The image is often pleasant to the eye but may lack details in certain area.

Author: John R. Victorine jvictor@kgs.ku.edu

References:

- (1) Color display of well logs, Peter L. Briggs, Mathematical Geology, Volume 17, Number 4, May 1985
- (2) Color Images of Kansas Subsurface Geology from Well Logs, D. R. Collins and J. H. Doveton, Computer & Geosciences, Vol. 12, No. 4B, pp.519-526 1986
- (3) Visualization of Subsurface Geology from Wireline Logs, David R. Collins, Digital Mapping Techniques '98-Workshop Proceedings U. S. Geological Survey Open-File Report 98-487
- (4) The Dakota Aquifer Program Annual Report, FY89 Kansas Geological Survey, Open-File Rept. 90-27 Annual Report, FY89-Appendix B
(http://www.kgs.ku.edu/Dakota/vol3/fy89/app_b.htm)
- (5) Composite Colour Display of Spectral Gamma-Ray Logs, by Matt Hall, Canadian Well Logging Society, Dec 2005, v.24
- (6) Geological Well Logs Their use in Reservoir Modeling, by Stefan M. Luthi, pg 84-86, ISBN 3-540-67840-9, © 2001 Springer-Verlag Berlin

The data displayed holds the initial information stored in the Log ASCII Standard (LAS) file and the Geologist Report ASCII Delimited file if the user loads the files from their PC. If the user loads the Log ASCII Standard (LAS) File from the KGS Server then the Well Header Information is automatically downloaded from the KGS Well Header Database Table.

Search KGS Database for Well Header Information Button

Displays a “Search for Data on KGS Server” Dialog that allows the user to .

Identification Panel

KGS & PC Primary KEY – Identification Numbers for the well

API-Number – API Number of Well
Well Status - Status, i.e. OIL, GAS, etc.

Name – Lease Name & Number

Other Well Information Panel

Operator Name & KGS Database KEY

Field Name & KGS Database KEY

Location Information Panel

State Name

County Name

Location

Township Range Section

XY Position

Latitude

Longitude

UTM Zone

UTM X Position

UTM Y Position

Z-Position

Depth – Total Depth of Well

Elevation – by Ground Level

Elevation – by Kelly Bushing

Elevation – by Derrick Floor

Comments – User Comments, not saved to the LAS version 3.0.

Header Information Source Buttons:

• **Show Initial Header Data** – Shows the Header Information initially loaded into Program.

• **Show KGS Well Header Data** – Shows the Header Information loaded from KGS Database.

Buttons:

Status – Displays “Select Status of Data” Dialog, user searches for the well status list for status of well.

Kansas TRS to Latitude, Longitude & Elevation – The buttons calls a KGS database routine to compute the Latitude, Longitude and Elevation from the Township, Range and Section.

Compute UTM – This button calls a UTM Java Math Package to convert Latitude & Longitude into UTM X, Y Coordinates.

OK – Transfer Data Values to Program

Close – Close this Dialog

NOTE: Initially the Basic Header information is loaded from the LAS version 2.0 file and other fields like Comments, Location are loaded from the Geologist Report Header Section.

The Header Information Dialog displays the contents of the header information data structure. The user can edit the fields and select the “Ok” Button to transfer the information back to the LAS File Viewer Program and any summary information will be updated in the LAS File Viewer Control and Plot.

As this example illustrates there are missing fields in the header information data. The user can select the “Search KGS Database for Well Header Information” Button, which will display a “Search for Data on KGS Server” Dialog that will allow the user to build a query that will download all wells that match the query.

This will display the “Search for Data on KGS Server” Dialog, see image below. This dialog allows the user to search the KGS database for well header data. In this example, the well of interest will be the Wellington KGS 1-32.

LTCI	API-Number	Well Name	Operator
	15-191-19025-...	WELLINGTON UNIT 58-INJ	TERRA RESOURCES,
LT	15-191-10272	DeTurk 3	Stelbar Oil Corp., Inc.
T	15-191-10054	WELLINGTON UNIT was Kamas 7 ...	Sinclair Prairie Oil Co.
T	15-191-10254	Wellington Unit 96	Stelbar Oil Corp. and D
T	15-191-43925	BARLOW 2	SHAWVER E B
T	15-191-19022	WELLINGTON UNIT - KAMAS LEAS...	COOPERATIVE REFGA
T	15-191-10296	Cora Stone 'A' 1	Stelbar Oil Corp., Inc.
LT	15-191-19021	Wellington Unit 141	Coop. Refining Assoc.
LT	15-191-22591	WELLINGTON KGS 1-32	BEREXCO LLC
T	15-191-10062	JOHN LUDWIG 1	STELBAR OIL CORP
T	15-191-43878	MURPHY 7	TRANSWESTERN OIL
T	15-191-10263	Wellington Unit 112	Stelbar Oil Corp., Inc.
T	15-191-10104	WELLINGTON UNIT, was PEASEL ...	SHAWVER E B
T	15-191-10100	WELLINGTON UNIT, was ERKER 9...	STELBAR OIL CORP IN

Search for Well Header Data in KGS Database Search By:

- **API-Number** – The user can search the KGS Database for well data by API-Number. The Format for the API is SS-CCC-99999 where
 - SS – Two Digit State Code
 - CCC – Three Digit County Code
 - 99999 – 5 Digit Well Number
- **Lease Name** – The user can search for well data by lease partial phrase, i.e. “Wellington”, which will look for all wells with the phrase “Wellington” in the lease name.
- **Township-Range-Section** – Search for a list of Wells by a specific area.

List of Kansas wells that match the search criteria

Load Well Header Buttons

- **Select** – Download the header information for the well selected.
- **Close** – Close this dialog

NOTE: LTCI Column in Table: L-LAS Files; T-Formation Tops; C-Measured Core Data; I-Core Images

As the Summary image suggests there are 3 methods for searching for the well header information within this dialog,

- By API-Number – KGS has a specific format for the API-Number, i.e.SS-CCC-99999 where SS is the state code for Kansas 15, CCC is the county code for Wellington KGS 1-32 it is 191 for Sumner County and the 5-Digit Well Number for the Wellington KGS 1-32 is 22591.

- By Partial Lease Name – The stored procedure used to retrieve the well header information allows the user to enter a partial phrase, in this example Wellington. The program places a ‘%’ in front and back of the phrase and sends the request to the Database, i.e. “%Wellington%”.

- By Township Range Section – This search is by location in Kansas, this search also allows the user to enter just the Township and Range to search for wells, e.g. to look for the Wellington KGS 1-32, enter Township as 31 set the S (South) Radio button and Range as 1 set the E (East) Radio button.

The user only needs to enter the above data and select the “Search” Button to display the list of Wells in the Kansas Database that match the search criteria. In the image below the Lease Name “Wellington” is entered to search for all wells in Kansas with the Phrase Wellington in it. The user searches through the list until they find the well of interest. In this example it is the Wellington KGS 1-32, which is highlighted.

LTCl	API-Number	Well Name	Operator
	15-191-19025-...	WELLINGTON UNIT 58-INJ	TERRA RESOURCES,
LT	15-191-10272	DeTurk 3	Stelbar Oil Corp., Inc.
T	15-191-10054	WELLINGTON UNIT was Kamas 7 ...	Sinclair Prairie Oil Co.
T	15-191-10254	Wellington Unit 96	Stelbar Oil Corp. and D
T	15-191-43925	BARLOW 2	SHAWVER E B
T	15-191-19022	WELLINGTON UNIT - KAMAS LEAS...	COOPERATIVE REFGA
T	15-191-10296	Cora Stone 'A' 1	Stelbar Oil Corp., Inc.
LT	15-191-19021	Wellington Unit 141	Coop. Refining Assoc.
LT_I	15-191-22591	WELLINGTON KGS 1-32	BEREXCO LLC
T	15-191-10062	JOHN LUDWIG 1	STELBAR OIL CORP
T	15-191-43878	MURPHY 7	TRANSWESTERN OIL
T	15-191-10263	Wellington Unit 112	Stelbar Oil Corp., Inc.
T	15-191-10104	WELLINGTON UNIT, was PEASEL ...	SHAWVER E B
T	15-191-10100	WELLINGTON UNIT, was ERKER 9...	STELBAR OIL CORP IN

The user clicks on the “Select” button to transfer the header information to the Edit Header Information Dialog.

LAS File Information

☒ Show Initial Header Data
☐ Show KGS Well Header Data

Search KGS Database for Well Header Information

Identification:

KGS Primary Key: PC Primary Key:

API-Number: Status:

Name:

Other Well Information:

Operator Name: Operator Code:

Field Name: Field Code:

Location Information:

State: County:

Location:

Township: Range: Section:

Kansas TRS to Latitude, Longitude & Elevation

XY Position:

Latitude: Longitude:

UTM Zone: Compute UTM

UTM-X: UTM-Y:

Z-Position:

Depth: Elevation: Kelly Bushing: Derrick Floor:

Comments:

Lease: Wellington KGS Well 1-32 (15-191-22591) ; operator: BEREXCO LLC; Field: Wellington
Location: T31S R1W, Sec. 32 : NE SW NE NE : 955 South, 877 West, from NE corner
Longitude: -97.4423481 ; Latitude: 37.3154639
County: Sumner
Total Depth: 3660 ; Elevation: 1259 GL

Ok Close

KGS Database Information

☐ Show Initial Header Data
☒ Show KGS Well Header Data

Search KGS Database for Well Header Information

Identification:

KGS Primary Key: PC Primary Key:

API-Number: Status:

Name:

Other Well Information:

Operator Name: Operator Code:

Field Name: Field Code:

Location Information:

State: County:

Location:

Township: Range: Section:

Kansas TRS to Latitude, Longitude & Elevation

XY Position:

Latitude: Longitude:

UTM Zone: Compute UTM

UTM-X: UTM-Y:

Z-Position:

Depth: Elevation: Kelly Bushing: Derrick Floor:

Comments:

Lat Long added from legal survey on intent. (DS-DRL 1-10-2011)

Ok Close

Location is from the Geologist report header section the 2nd line of the ASCII Text file.

Comments are from the Geologist report header section the lines before the start of the data in the ASCII Text file.

The “Show KGS Well Header Data” radio button will become enabled if the KGS well header information has been downloaded. The user can move between the initial header information by selecting the “Show Initial Header Data” radio button the KGS well header information by selecting the “Show KGS Well Header Data” radio button.

There a number of buttons on the panel that will allow the user to change or compute data in the header information. The “Kansas TRS to Latitude, Longitude & Elevation” Button computes the latitude, longitude and elevation from the township, range and section by making an Oracle PL/SQL Stored Procedure, e.g.

http://chasm.kgs.ku.edu/ords/iqstrat.TRS2LL_pkg.getXML?iTownship=31&sTownship=S&iRange=1&sRange=E&iSection=32

The call will return a XML with the latitude, longitude and ground level elevation.

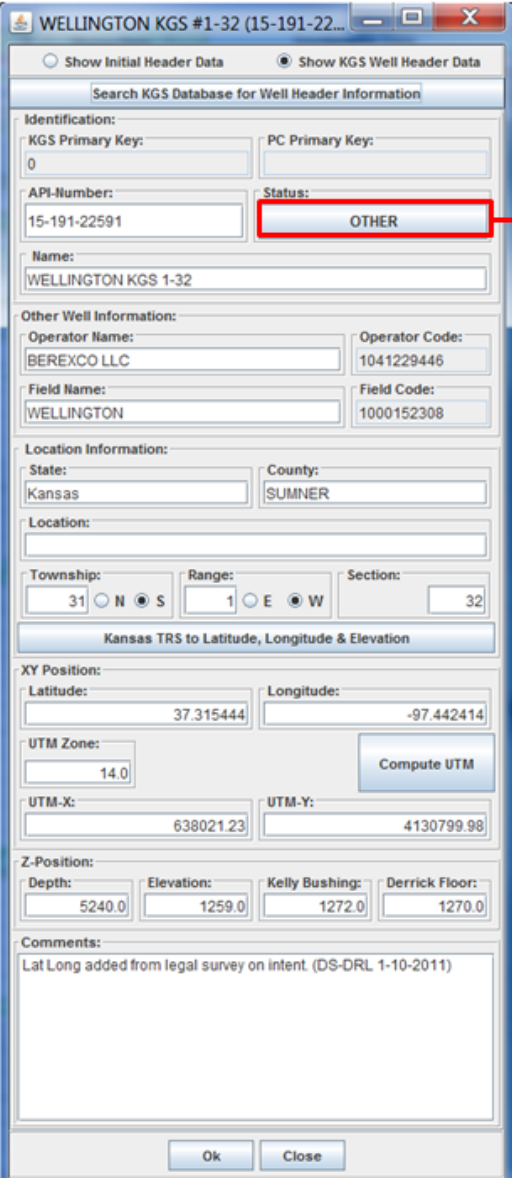
```

<?xml version="1.0"?>
<!DOCTYPE headers [
<!ELEMENT headers (data*)>
<!ATTLIST headers records CDATA #IMPLIED>
<!ELEMENT data (info*,
                other?,
                loc?,
                xy?,
                z?,
                comments?,
                misc?,
                cnt?)>
<!ELEMENT info EMPTY>
<!ATTLIST info kid CDATA #IMPLIED
               well_kid CDATA #IMPLIED
               key CDATA #IMPLIED
               type CDATA #IMPLIED
               api CDATA #IMPLIED
               name CDATA #IMPLIED
               status CDATA #IMPLIED
               error CDATA #IMPLIED>
<!ELEMENT other EMPTY>
<!ATTLIST other operator CDATA #IMPLIED
               oper_kid CDATA #IMPLIED
               field CDATA #IMPLIED
               field_kid CDATA #IMPLIED>
<!ELEMENT loc EMPTY>
<!ATTLIST loc state CDATA #IMPLIED
               state_cd CDATA #IMPLIED
               county CDATA #IMPLIED
               county_cd CDATA #IMPLIED
               loc CDATA #IMPLIED
               town CDATA #IMPLIED
               town_dir CDATA #IMPLIED
               range CDATA #IMPLIED
               range_dir CDATA #IMPLIED
               section CDATA #IMPLIED>
<!ELEMENT xy EMPTY>
<!ATTLIST xy latitude CDATA #IMPLIED
               longitude CDATA #IMPLIED
               zone CDATA #IMPLIED
               utm_x CDATA #IMPLIED
               utm_y CDATA #IMPLIED>
<!ELEMENT z EMPTY>
<!ATTLIST z depth CDATA #IMPLIED
               gl CDATA #IMPLIED
               kb CDATA #IMPLIED
               df CDATA #IMPLIED>
<!ELEMENT comments (#PCDATA)>
<!ELEMENT misc EMPTY>
<!ATTLIST misc user CDATA #IMPLIED
               access CDATA #IMPLIED
               source CDATA #IMPLIED
               date CDATA #IMPLIED>
<!ELEMENT cnt EMPTY>
<!ATTLIST cnt las CDATA #IMPLIED
               tops CDATA #IMPLIED
               core CDATA #IMPLIED
               images CDATA #IMPLIED]>
<headers records="1">
  <data>
    <loc town="31" town_dir="S" range="1" range_dir="E" section="32" />
    <xy latitude="37.311703" longitude="-97.339619" />
    <z gl="1277" />
  </data>
</headers>

```

The “UTM” Button will compute the UTM XY coordinates from the latitude & longitude. The analysis uses A Working Manual by John P. Snyder, U.S. Geological Survey Professional Paper 1395, USG Printing Office, Washington, DC, 1987 (http://pubs.er.usgs.gov/djvu/PP/PP_1395.pdf).

The Status button allows the user to change the well status.



The 'WELLINGTON KGS #1-32 (15-191-22...)' window displays various well header information. The 'Status' button is highlighted in red. A red arrow points from this button to the 'Select Status of Data' dialog box.

The 'Select Status of Data' dialog box displays a list of possible common well status options. The 'LOC' status is highlighted in yellow.

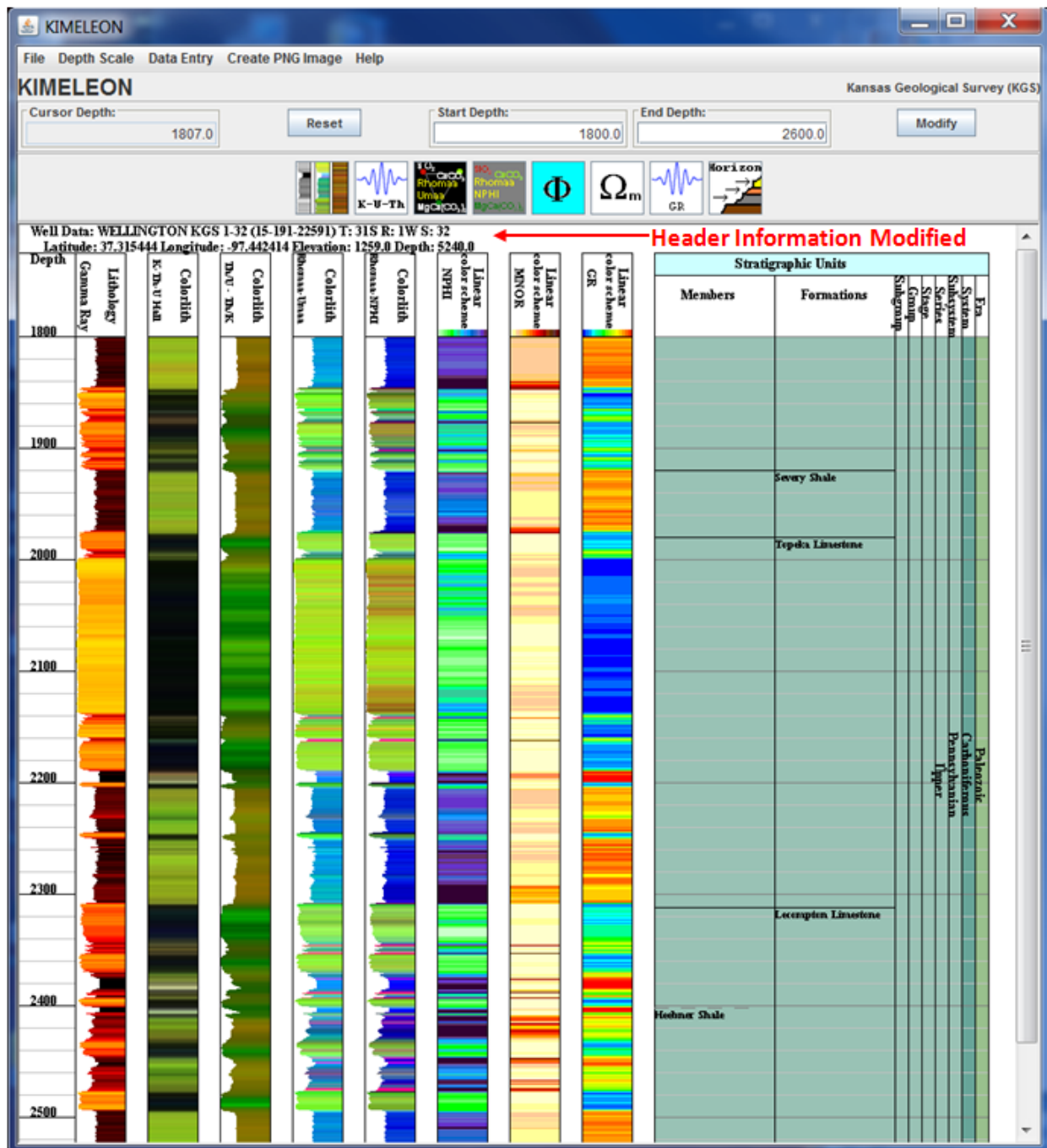
Mnemonic	Description
LOC	Location
INTENT	Intent
D&A	Dry & Abandoned
SUS	Suspended
P&A	Plugged & Abandoned
EOR	Enhanced Oil Recovery
EOR-P&A	Enhanced Oil Recovery - Plugged & Abandoned
OIL	Oil
D&A-O	Dry & Abandoned - Oil Show
SUS-O	Suspended - Oil
OIL-P&A	Oil - Plugged & Abandoned
GAS	Gas
D&A-G	Dry & Abandoned - Gas Show
SUS-G	Suspended - Gas
GAS-P&A	Gas - Plugged & Abandoned
O&G	Oil & Gas
D&A-OG	Dry & Abandoned - Oil & Gas Show
SUS-OG	Suspended - Oil & Gas
O&G-P&A	Oil & Gas - Plugged & Abandoned
OTHER	Other
OTHER-P&A	Other - Plugged & Abandoned
INJ	Injection
INJ-P&A	Injection - Plugged & Abandoned
SWD	Salt Water Disposal
SWD-P&A	Salt Water Disposal - Plugged & Abandoned
CBM	Coal Bed Methane
CBM-P&A	Coal Bed Methane - Plugged & Abandoned
OUTCROP	Measured Section

The dialog box includes 'Select' and 'Cancel' buttons at the bottom.

Select “Status” Button to display the “Select Status of Data” Dialog, which displays the list of possible common well status.

Highlight the desired status and click on the “Select” Button to transfer the Mnemonic to the “Status” Button text.

Select the “Ok” Button to update the Header Information in plot panel.



Adding & Modifying Tops

Enter Horizon Data - Stratigraphic Units Panel

The Stratigraphic Units Panel allows the user to Add, Modify or Remove Stratigraphic Units. There are two files that are used to assist in mapping Stratigraphic Units to a specific top name.

- The 2010 International Commission on Stratigraphy Stratigraphic Units and RGB color for the Stratigraphic Units XML File (<http://www.kgs.ku.edu/software/gemini/data/ISC.xml>) are used to display the tops in the Stratigraphics Plot Track by Age (RGB Color).
- The Kansas Geological Survey (KGS) Stratigraphic Succession in Kansas, edited by D.E. Zeller, December 1968, updated 2012 (<http://www.kgs.ku.edu/software/gemini/data/kansas.xml>) will help map the Kansas Top Names to Stratigraphic Units, System, Sub-System, Series, etc. and to map the top depth of one top pick to the base depth of another top pick.

Starting Depth & Ending Depth of Stratigraphic Name

Confidence Level of the tops selection.

Stratigraphic Unit Rank radio buttons, defines & sets the location of the unit on the Stratigraphic Unit Plot Track.

Stratigraphic Name & Alternate Name

Add/Modify – Move data to Table.
Clear – Clear all text fields.

List of Stratigraphic Units (Tops).

International Commission on Stratigraphic Units. User selects the ICS Chart button to display Standard Units.

Stratigraphic Name belongs to section. Allows the user to place a member, bed, etc. with a formation, group, etc.

“Stratigraphic Units Selected” Table.

Table Buttons
Modify – Modify the Stratigraphic Units Data.
Remove – Remove Stratigraphic Unit from table.
Remove All – Clear all content Units from the table.

The KGS Stratigraphic Succession of Kansas edited by D. E. Zeller, Dec. 1968, updated 2012 (http://www.kgs.ku.edu/PRS/Ozark/TYPE_LOG/Stratigraphic/index.html).

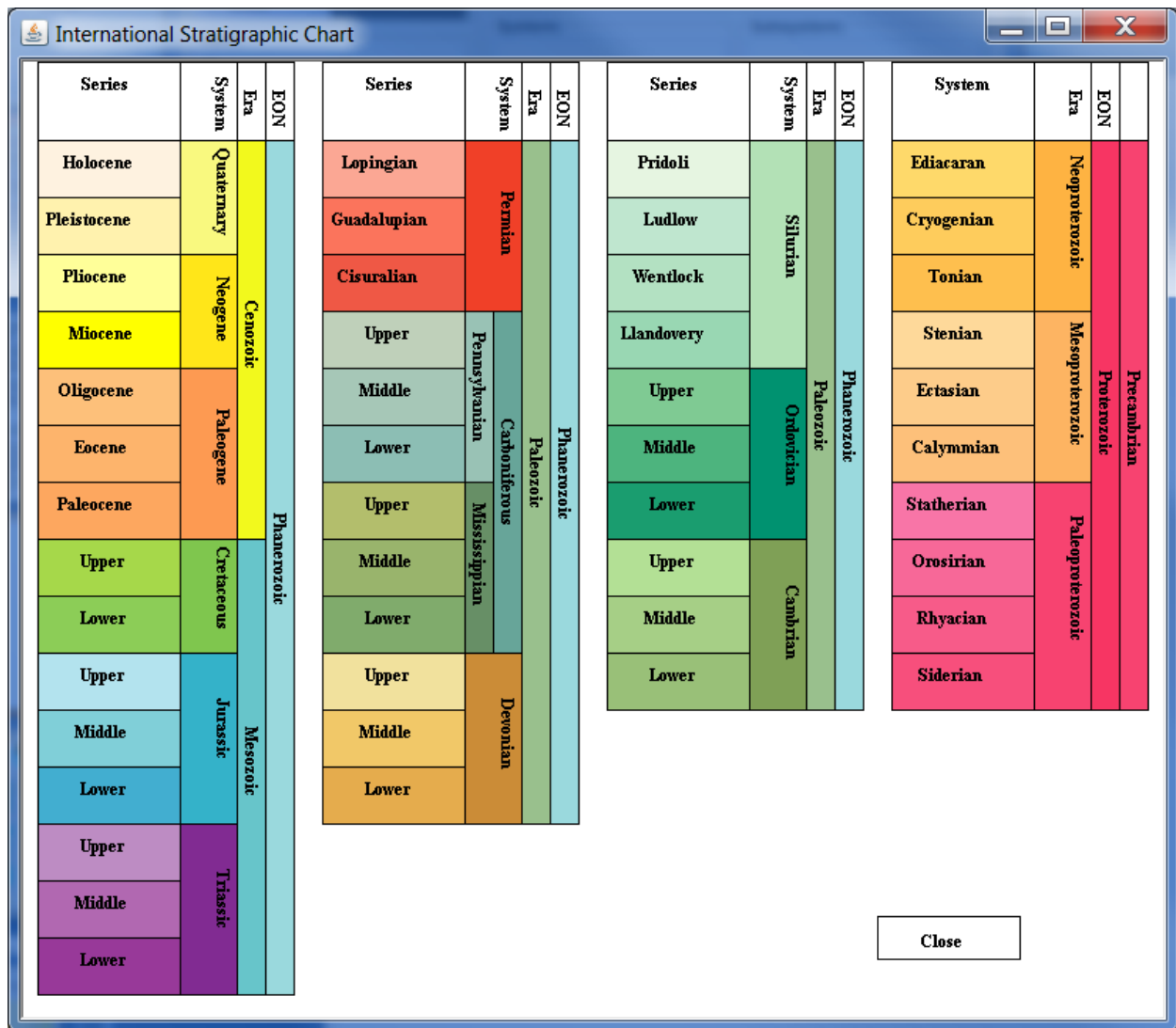
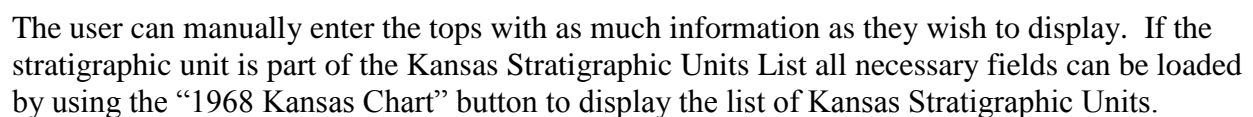


Figure: The 2010 International Commission on Stratigraphy Stratigraphic Units and RGB colors for the Stratigraphic Units.

This first example is to add a Shawnee Group, which is part of the KGS Stratigraphic Succession in Kansas. Select the “Data Entry” Menu and then click on the “Data Entry - Formation Tops” menu option to display the “Enter Horizon Data” Dialog with the “Stratigraphic Units” dialog.



The Topeka Limestone, Lecompton Limestone and the Heebner Shale belong to the Shawnee Group, set the Starting Depth text field to 1980 and the Ending Depth text field to 2402. Click on the “1968 Kansas Chart” Button to display the State Stratigraphic Units Dialog.

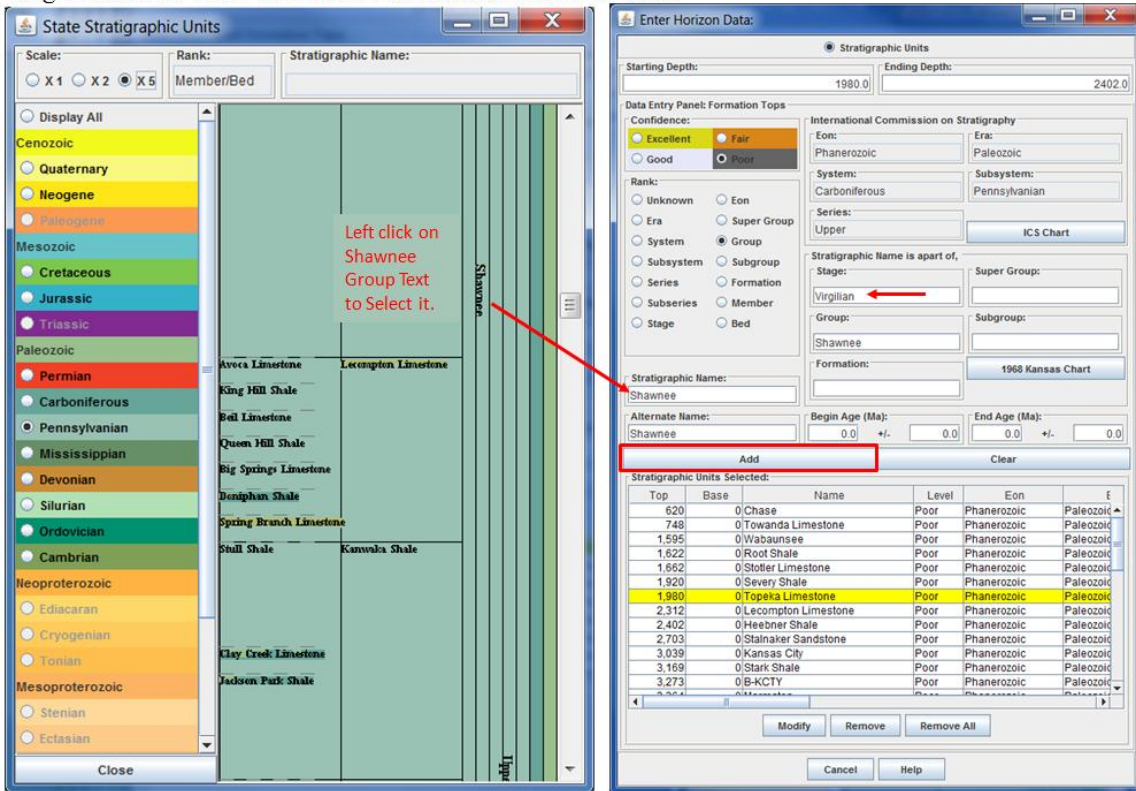
The left screenshot shows the 'Enter Horizon Data' window. The 'Starting Depth' is 1980.0 and the 'Ending Depth' is 2402.0. The '1968 Kansas Chart' button is highlighted with a red box. The right screenshot shows the 'State Stratigraphic Units' dialog. The 'Scale' is X1 and the 'Rank' is Member/Bed. The 'Stratigraphic Name' field is empty. The 'Display All' radio button is selected. The 'Stratigraphic Units' list on the right shows the 'Shawnee' group highlighted in yellow.

Change the Scale to X5, select the “Pennsylvanian” radio button.

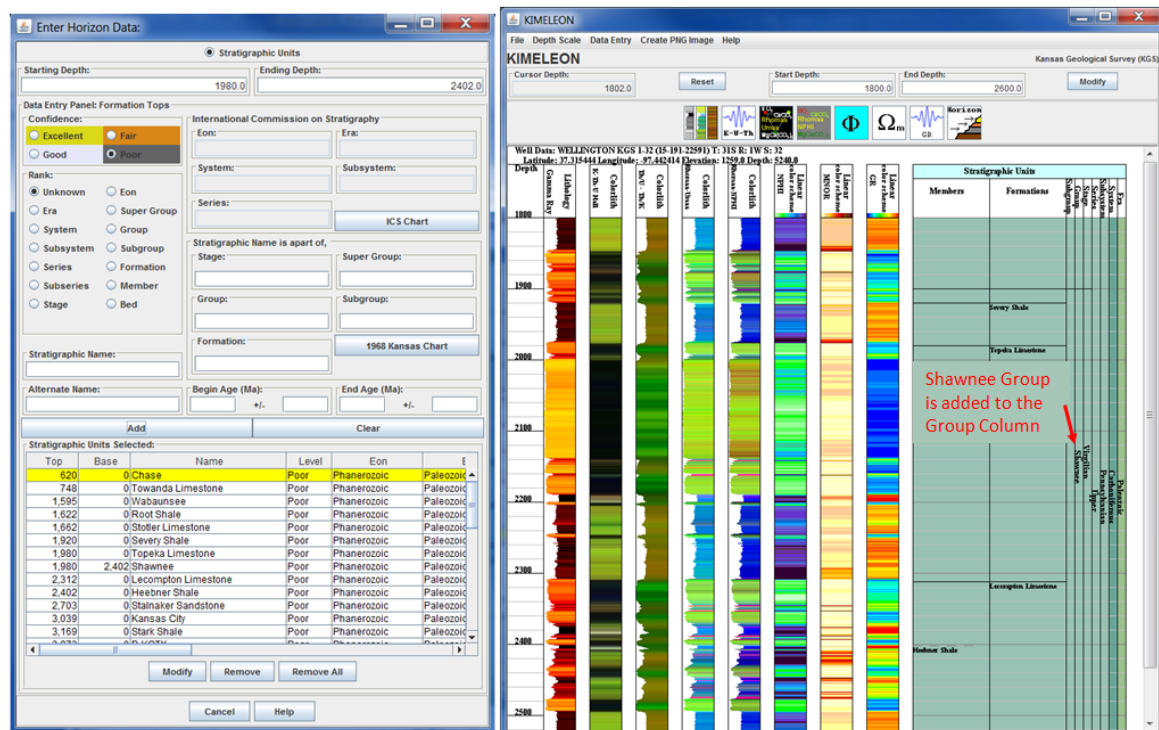
Scroll down to the Lecompton Limestone Formation. Notice it is part of the Shawnee Group.

The left screenshot shows the 'State Stratigraphic Units' dialog with the 'Scale' set to X5 and the 'Pennsylvanian' radio button selected. The right screenshot shows the 'Lecompton Limestone' formation highlighted in yellow, with a red box around the 'Shawnee' group label.

Left mouse click on the Shawnee Group text to transfer all the Stratigraphic Units to the “Stratigraphic Units” Panel on the “Enter Horizon Data” Dialog. Notice also that the Shawnee Group belongs to the Virgilian Stage. Select the “Add” Button to move Shawnee into the table.



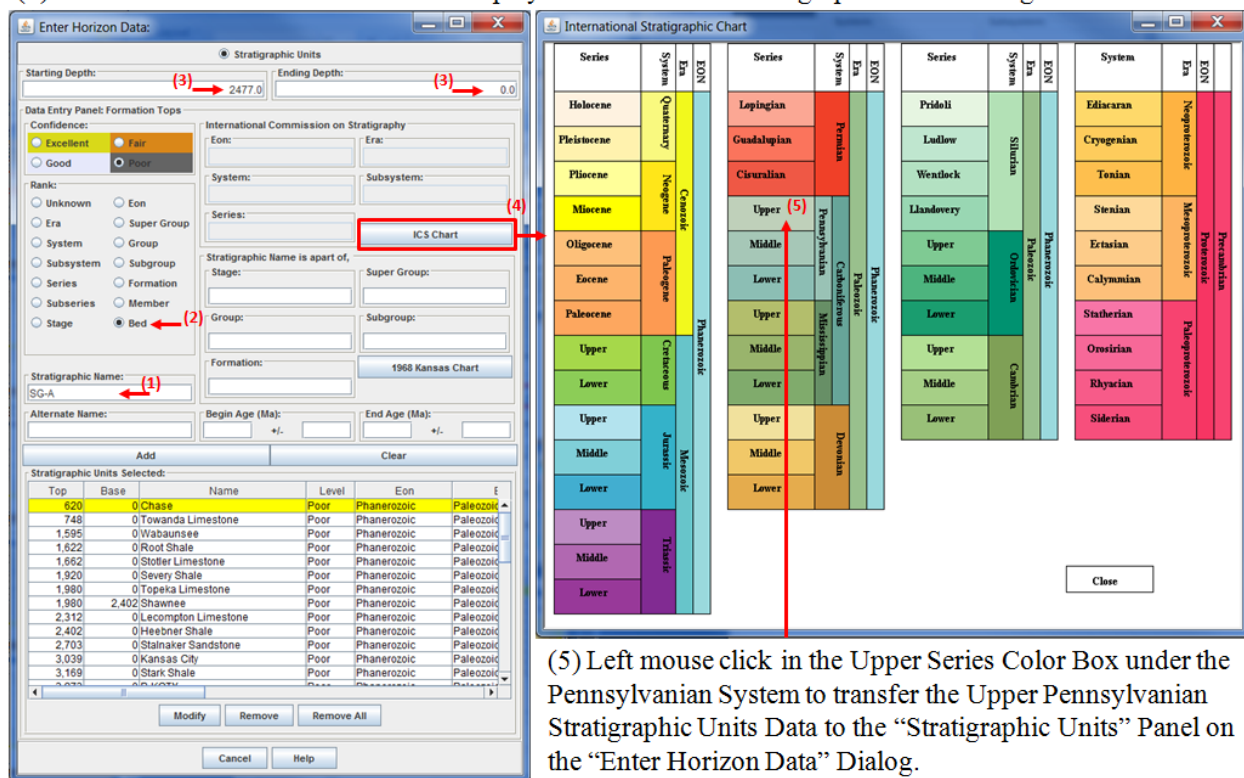
Notice that the Shawnee Group is added to the Group Column of the Stratigraphic Units Plot Track. Shawnee was added to the “Stratigraphic Units Selected” table, see the highlighted top in the table.



Add Unknown Bed (SG-A Bed) to the Stratigraphic Units List

This second example is to add an unknown bed (SG-A Bed) to the Stratigraphic Units List, which is not part of the KGS Stratigraphic Succession in Kansas. This example will enter the basic information to the Depth Range text fields, the Stratigraphic Name text field and setting the Rank as a BED. The Stratigraphic Units will be provided by the International Commission on Stratigraphy Dialog by selecting the “ICS Chart” Button. The “International Stratigraphic Chart” Dialog displays the stratigraphic units as a series of colored boxes, each stratigraphic unit is has the recommended RGB (Red-Green-Blue) Color defined by the International Commission on Stratigraphy. The user only needs to click the colored box to select all the stratigraphic data associated with the selected stratigraphic unit and transfer the data back to the “Stratigraphic Units” Panel in the “Enter Horizon Data” Dialog.

- (1) Add SG-A to the “Stratigraphic Name” Text field.
- (2) Select the Bed Radio Button in the Rank Panel
- (3) Set the Starting Depth to 2477.0 and the Ending Depth to 0.0 Limestone Depth Range.
- (4) Click on the “ICS Chart” Button to display the International Stratigraphic Chart Dialog.



- (5) Left mouse click in the Upper Series Color Box under the Pennsylvanian System to transfer the Upper Pennsylvanian Stratigraphic Units Data to the “Stratigraphic Units” Panel on the “Enter Horizon Data” Dialog.

Enter Horizon Data:

Stratigraphic Units

Starting Depth: 2477.0 Ending Depth: 0.0

Data Entry Panel: Formation Tops

Confidence: ☒ Excellent ☐ Fair ☐ Good ☐ Poor

Rank: ☐ Unknown ☐ Eon ☐ Era ☐ Super Group ☐ System ☐ Group ☐ Subsystem ☐ Subgroup ☐ Series ☐ Formation ☐ Subseries ☐ Member ☐ Stage ☒ Bed

Stratigraphic Name: SG-A

Alternate Name: SG-A

Begin Age (Ma): +/- End Age (Ma): +/-

Add Clear

Stratigraphic Units Selected:

Top	Base	Name	Level	Eon	E
620	0	Chase	Poor	Phanerozoic	Paleozoic
748	0	Towanda Limestone	Poor	Phanerozoic	Paleozoic
1,595	0	Wabaunsee	Poor	Phanerozoic	Paleozoic
1,622	0	Root Shale	Poor	Phanerozoic	Paleozoic
1,662	0	Stotter Limestone	Poor	Phanerozoic	Paleozoic
1,920	0	Severy Shale	Poor	Phanerozoic	Paleozoic
1,980	0	Topeka Limestone	Poor	Phanerozoic	Paleozoic
1,980	2,402	Shawnee	Poor	Phanerozoic	Paleozoic
2,312	0	Lecompton Limestone	Poor	Phanerozoic	Paleozoic
2,402	0	Heebner Shale	Poor	Phanerozoic	Paleozoic
2,703	0	Stalnaker Sandstone	Poor	Phanerozoic	Paleozoic
3,039	0	Kansas City	Poor	Phanerozoic	Paleozoic
3,169	0	Stark Shale	Poor	Phanerozoic	Paleozoic

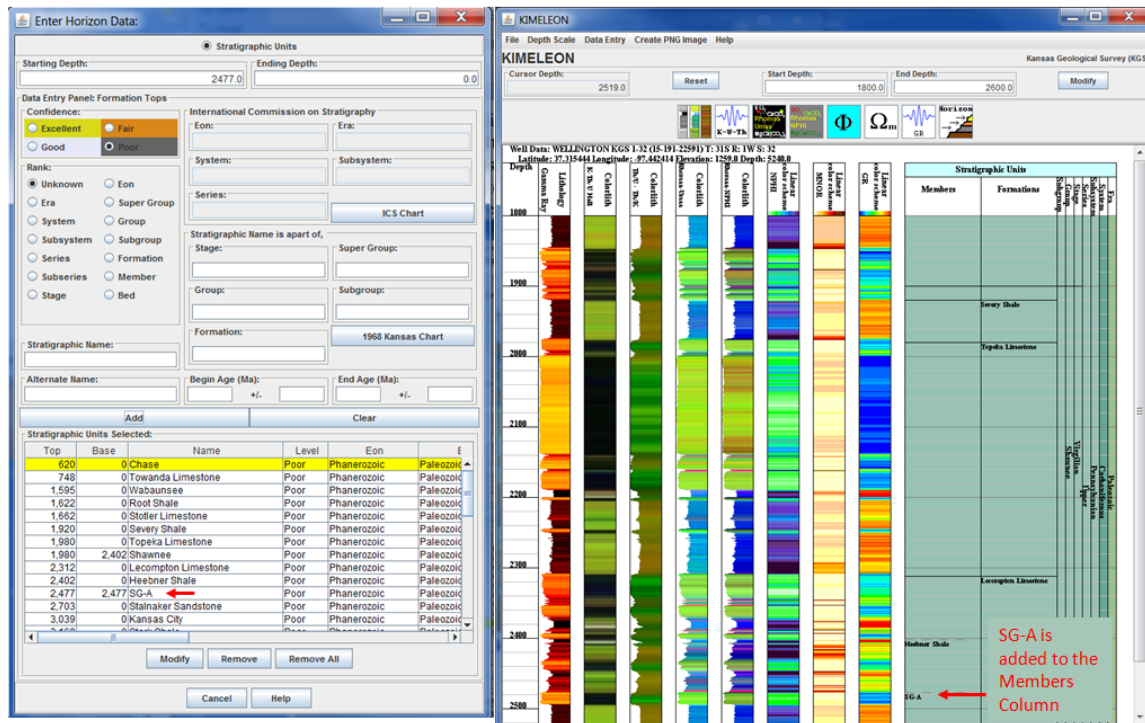
Modify Remove Remove All

Cancel Help

Notice that the Stratigraphic Units Data from the ICS Chart is transferred to the "International Commission on Stratigraphy" Panel.

Now Select the "Add" Button to transfer the Stratigraphic Units Data for the SG-A Bed to the "Stratigraphic Units Selected" Table.

Notice that the SG-A Bed is added to the Members Column of the Stratigraphic Units Plot Track. SG-A Bed was added to the "Stratigraphic Units Selected" table, see the highlighted top in the table.



Modify data for the Severy Shale Formation in the Stratigraphic Units List

Highlight the Severy Shale, select the “Modify” Button.

Enter Horizon Data: Stratigraphic Units

Starting Depth: 1920.0 Ending Depth: 0.0

Confidence: ☒ Excellent ☐ Fair ☐ Good ☐ Poor

International Commission on Stratigraphy

Eon: Era:

System: Subsystem:

Series: ICS Chart

Stratigraphic Name is apart of, Stage: Super Group:

Group: Subgroup:

Formation: 1968 Kansas Chart

Stratigraphic Name:

Alternate Name: Begin Age (Ma): End Age (Ma):

Add Clear

Stratigraphic Units Selected:

Top	Base	Name	Level	Eon	f
620	0	Chase	Poor	Phanerozoic	Paleozoic
748	0	Towanda Limestone	Poor	Phanerozoic	Paleozoic
1595	0	Wabunsee	Poor	Phanerozoic	Paleozoic
1622	0	Root Shale	Poor	Phanerozoic	Paleozoic
1662	0	Stoller Limestone	Poor	Phanerozoic	Paleozoic
1920	0	Severy Shale	Poor	Phanerozoic	Paleozoic
1980	0	Topeka Limestone	Poor	Phanerozoic	Paleozoic
1980	2,402	Shawnee	Poor	Phanerozoic	Paleozoic
2,312	0	Lecompton Limestone	Poor	Phanerozoic	Paleozoic
2,402	0	Heebner Shale	Poor	Phanerozoic	Paleozoic
2,477	2,477	SG-A	Poor	Phanerozoic	Paleozoic
2,703	0	Stalnaker Sandstone	Poor	Phanerozoic	Paleozoic
3,039	0	Kansas City	Poor	Phanerozoic	Paleozoic

Modify Remove Remove All

Cancel Help

Notice the Stratigraphic Units Data are loaded into the Stratigraphic Units Panel Fields and the “Stratigraphic Name is apart of,” fields are empty.

Enter Horizon Data: Stratigraphic Units

Starting Depth: 1920.0 Ending Depth: 0.0

Confidence: ☒ Excellent ☐ Fair ☐ Good ☐ Poor

International Commission on Stratigraphy

Eon: Era:

System: Subsystem:

Series: ICS Chart

Stratigraphic Name is apart of, Stage: Super Group:

Group: Subgroup:

Formation: 1968 Kansas Chart

Stratigraphic Name:

Alternate Name: Begin Age (Ma): End Age (Ma):

Modify Clear

Stratigraphic Units Selected:

Top	Base	Name	Level	Eon	f
620	0	Chase	Poor	Phanerozoic	Paleozoic
748	0	Towanda Limestone	Poor	Phanerozoic	Paleozoic
1595	0	Wabunsee	Poor	Phanerozoic	Paleozoic
1622	0	Root Shale	Poor	Phanerozoic	Paleozoic
1662	0	Stoller Limestone	Poor	Phanerozoic	Paleozoic
1920	0	Severy Shale	Poor	Phanerozoic	Paleozoic
1980	0	Topeka Limestone	Poor	Phanerozoic	Paleozoic
1980	2,402	Shawnee	Poor	Phanerozoic	Paleozoic
2,312	0	Lecompton Limestone	Poor	Phanerozoic	Paleozoic
2,402	0	Heebner Shale	Poor	Phanerozoic	Paleozoic
2,477	2,477	SG-A	Poor	Phanerozoic	Paleozoic
2,703	0	Stalnaker Sandstone	Poor	Phanerozoic	Paleozoic
3,039	0	Kansas City	Poor	Phanerozoic	Paleozoic

Modify Remove Remove All

Cancel Help

Left mouse click on the Severy Shale text to transfer all the Stratigraphic Units to the “Stratigraphic Units” Panel on the “Enter Horizon Data” Dialog. Notice also that the Severy Shale belongs to the Wabunsee.

State Stratigraphic Units

Scale: ☐ X1 ☐ X2 ☒ X5

Rank: Member/Bed

Stratigraphic Name:

Display All

Cenozoic

- ☐ Quaternary
- ☐ Neogene
- ☐ Paleogene

Mesozoic

- ☐ Cretaceous
- ☐ Jurassic
- ☐ Triassic

Paleozoic

- ☒ Permian
- ☐ Carboniferous
- ☐ Pennsylvanian
- ☐ Mississippian
- ☐ Devonian
- ☐ Silurian
- ☐ Ordovician
- ☐ Cambrian

Neoproterozoic

- ☐ Ediacaran
- ☐ Cryogenian
- ☐ Tonian

Mesoproterozoic

- ☐ Stenian
- ☐ Ectasian
- ☐ Calymnian

Paleoproterozoic

- ☐ Gaskiers
- ☐ Sturtian
- ☐ Dryas
- ☐ Marinoan

Close

Left click on Severy Shale Text to Select it.

Enter Horizon Data: Stratigraphic Units

Starting Depth: 1920.0 Ending Depth: 1980.0

Confidence: ☒ Excellent ☐ Fair ☐ Good ☐ Poor

International Commission on Stratigraphy

Eon: Era:

System: Subsystem:

Series: ICS Chart

Stratigraphic Name is apart of, Stage: Super Group:

Group: Subgroup:

Formation: 1968 Kansas Chart

Stratigraphic Name:

Alternate Name: Begin Age (Ma): End Age (Ma):

Modify Clear

Stratigraphic Units Selected:

Top	Base	Name	Level	Eon	f
620	0	Chase	Poor	Phanerozoic	Paleozoic
748	0	Towanda Limestone	Poor	Phanerozoic	Paleozoic
1595	0	Wabunsee	Poor	Phanerozoic	Paleozoic
1622	0	Root Shale	Poor	Phanerozoic	Paleozoic
1662	0	Stoller Limestone	Poor	Phanerozoic	Paleozoic
1920	0	Severy Shale	Poor	Phanerozoic	Paleozoic
1980	0	Topeka Limestone	Poor	Phanerozoic	Paleozoic
1980	2,402	Shawnee	Poor	Phanerozoic	Paleozoic
2,312	0	Lecompton Limestone	Poor	Phanerozoic	Paleozoic
2,402	0	Heebner Shale	Poor	Phanerozoic	Paleozoic
2,477	2,477	SG-A	Poor	Phanerozoic	Paleozoic
2,703	0	Stalnaker Sandstone	Poor	Phanerozoic	Paleozoic
3,039	0	Kansas City	Poor	Phanerozoic	Paleozoic

Modify Remove Remove All

Cancel Help

Enter Horizon Data:

Stratigraphic Units

Starting Depth: 1920.0 Ending Depth: 1980.0

Data Entry Panel: Formation Tops

Confidence: ☒ Excellent ☐ Fair ☐ Good ☐ Poor

Rank: ☐ Unknown ☐ Eon ☐ Era ☐ Super Group ☐ System ☐ Group ☐ Subsystem ☐ Subgroup ☒ Series ☐ Formation ☐ Subseries ☐ Member ☐ Stage ☐ Bed

Stratigraphic Name: Severy Shale

Alternate Name: Severy Shale

Begin Age (Ma): 0.0 +/- End Age (Ma): 0.0 +/-

Modify Clear

Stratigraphic Units Selected:

Top	Base	Name	Level	Eon	System	Subsystem	Series	Subgroup
620	0	Chase	Poor	Phanerozoic	Paleozoic			
748	0	Towanda Limestone	Poor	Phanerozoic	Paleozoic			
1,595	0	Wabunsee	Poor	Phanerozoic	Paleozoic			
1,622	0	Root Shale	Poor	Phanerozoic	Paleozoic			
1,662	0	Stoller Limestone	Poor	Phanerozoic	Paleozoic			
1,920	0	Severy Shale	Poor	Phanerozoic	Paleozoic			
1,980	0	Topeka Limestone	Poor	Phanerozoic	Paleozoic			
1,980	2,402	Shawnee	Poor	Phanerozoic	Paleozoic			
2,312	0	Lecompton Limestone	Poor	Phanerozoic	Paleozoic			
2,402	0	Heebner Shale	Poor	Phanerozoic	Paleozoic			
2,477	2,477	SG-A	Poor	Phanerozoic	Paleozoic			
2,703	0	Stalaker Sandstone	Poor	Phanerozoic	Paleozoic			
3,039	0	Kansas City	Poor	Phanerozoic	Paleozoic			

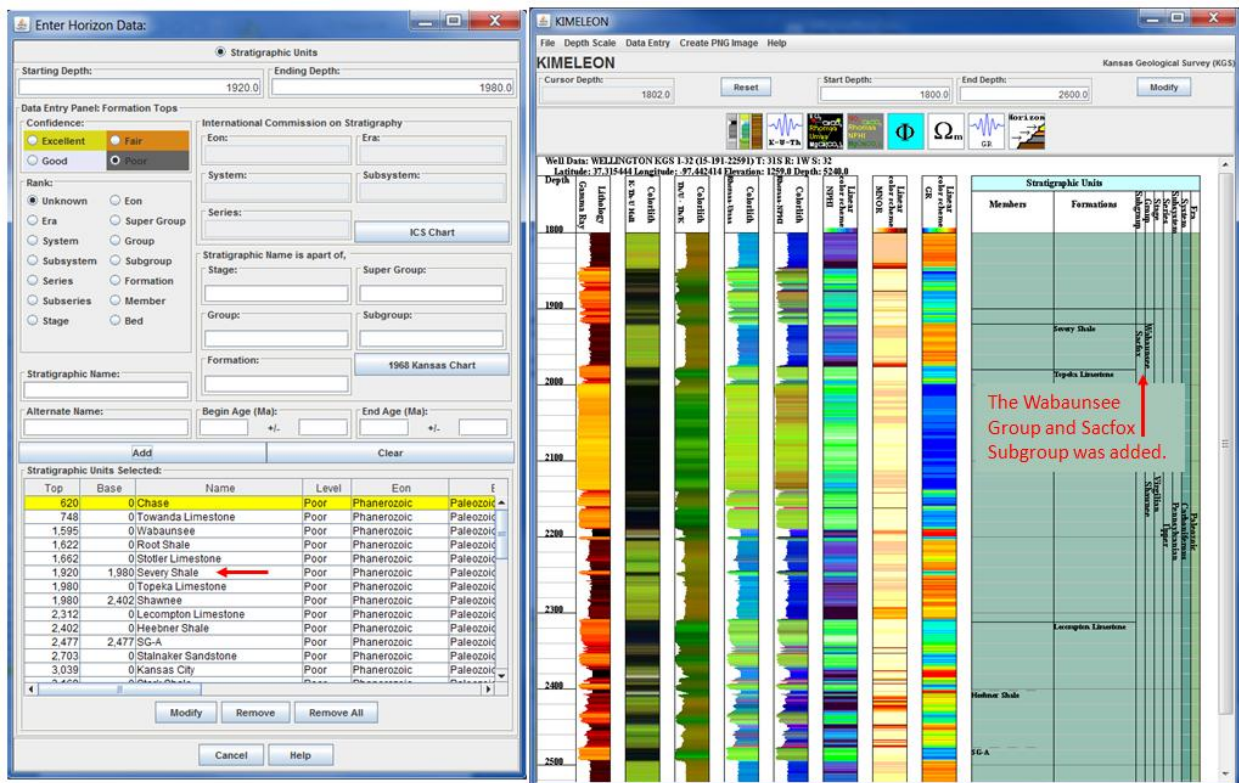
Modify Remove Remove All

Cancel Help

Change the Ending Depth to 1980.0, which is the start Depth of the Topeka Limestone and the next Formation.

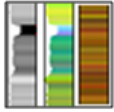
Now Select the "Modify" Button to save the changes and modify the contents of the "Severy Shale" Stratigraphic Units in the "Stratigraphic Units Selected" table.

Notice that the Wabunsee Group & the Sacfox Subgroup were added.



KIMELEON Colorlith & Color Image Tracks

The KIMELEON Plot Panels allows the user to modify the colorlith & color image plot tracks.

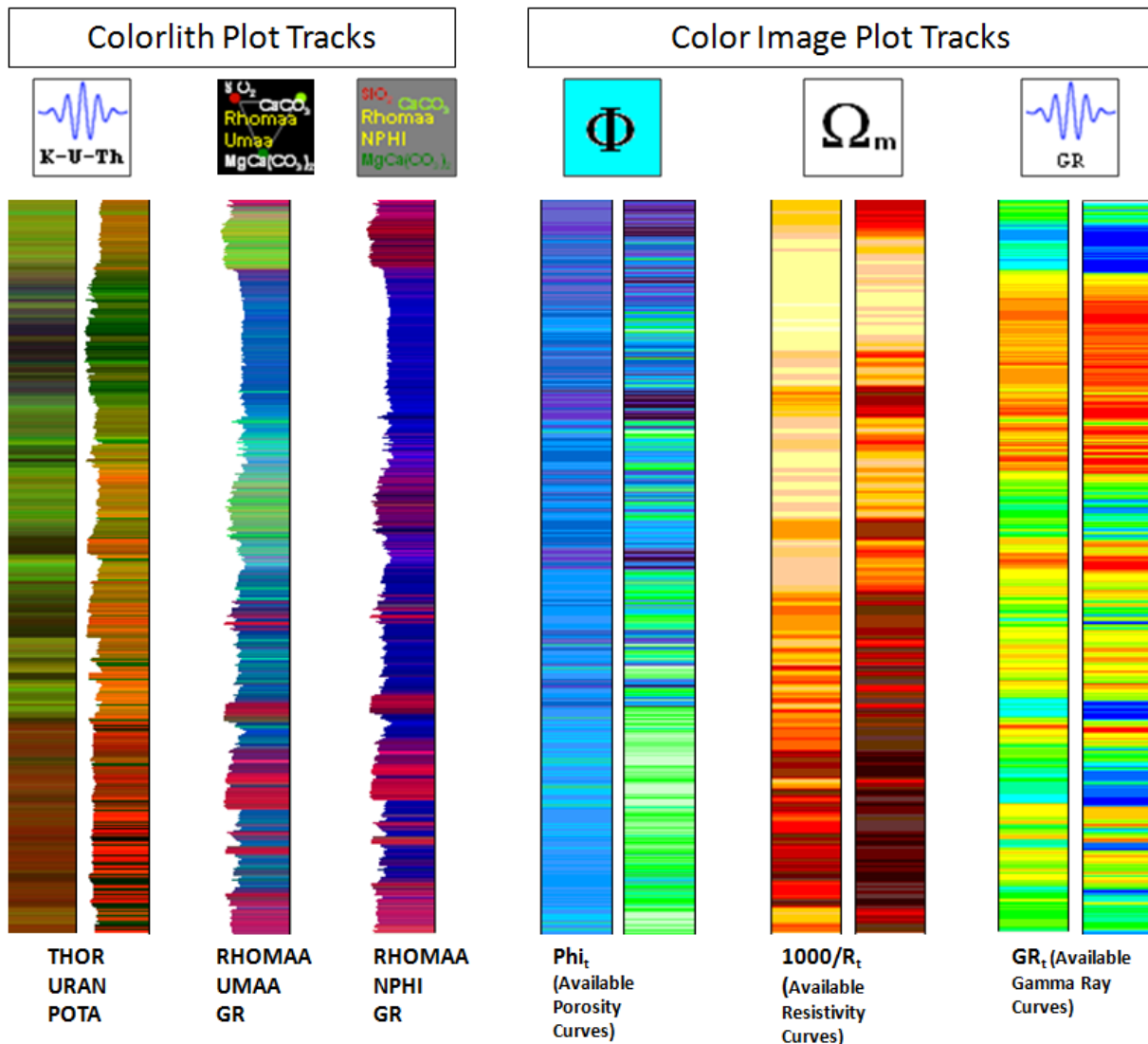


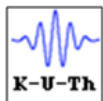
Summary of all image Colorlith & Color Image Plot Tracks



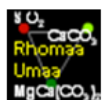
Summary of all image Colorlith & Color Image Plot Tracks against formation tops picks.

There are two types of Image Tracks, the Colorlith, which maps 3 log curves to the each Red, Green and Blue color and the Color Image Tracks that uses up to 16 individual colors to plot a single log curve values.





Spectral Gamma Ray Image Panel – Thorium (Th), Uranium (U) and Potassium (K) Log Curves and the K-Th-U Hall Colorlith and Spectral Gamma Ray Ratio Colorlith Plot Tracks.



Rhomaa-Umaa Image Panel – Apparent Grain Density (RHOMaa), Apparent Photoelectric (Umaa) and Gamma Ray (GR) Log Curves and the Rhomaa-Umaa Colorlith Plot Tracks.



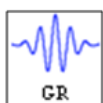
Rhomaa-NPHI Image Panel – Apparent Grain Density (RHOMaa), Neutron Porosity (NPHI) and Gamma Ray (GR) Log Curves and the Rhomaa-NPHI Colorlith Plot Tracks.



Porosity Image Panel – Any Available Porosity Log Curve Type and the Linear & Non-Linear Color Image Plot Tracks.



Conductivity Image Panel – Any Available Resistivity Log Curve Type and the Linear & Non-Linear Color Image Plot Tracks.



Gamma Ray Image Panel – Any Available Gamma Ray Log Curve Type and the Linear & Non-Linear Color Image Plot Tracks.

Colorlith Plot Tracks

The Colorlith format uses 3 log curves and maps each log curve to a red, green or blue color. Each color has a range from 0 to 255, where 0 the darkest color value (“black”) and 255 is lightest color value (“white”). The initial minimum & maximum log curve values have been set by previous authors^(2,3,4,5), which initially suggests lithology by color. This program was designed to over write the color limits by log curve minimum and maximum to change the basic presentation of the RGB track to bring out subtle color changes by depth. There are three sets of log curves that are presented in the colorlith format, which are fixed curves and cannot be altered. The log curves types are as follows,

- Spectral Gamma Ray Log Curves K-Th-U Hall Colorlith Plot Track, which uses all the spectral gamma ray log curves to create the RGB color schema, which was suggested by Matt Hall⁽⁵⁾ to map the “energy” levels of the gamma-ray emissions to the visible spectrum.
 - POTA – Potassium Concentration (potassium-40 nuclei having the lowest energy) assigned a red color range initially from 0 to 5 %.
 - THOR – Thorium Concentration (thorium nuclei having the next highest energy) assigned a green color range initially from 0 to 20 ppm.
 - URAN – Uranium Concentration (uranium nuclei having the highest energy) assigned a blue color range initially from 0 to 30 ppm.
- The Spectral Gamma Ray Ratio Curves Th/U and Th/K Colorlith Plot Track, which maps each ratio to a specific color scheme suggested by Doveton⁽⁴⁾.
 - Th/U – Thorium/Uranium Ratio assigned a red color range initially from 0.57 to 24.5.

- Th/K – Thorium/Potassium Ratio assigned a green color range initially from 1.02 to 24.5.
- The Litho-Density Log Curves with Photoelectric Factor (PE) Colorlith Plot Track, which uses the Bulk Density (RHOB), Neutron Porosity (NPHI) and the Photoelectric Factor (PE) log curves to compute the RHOMaa-Umaa curves, the limits suggested by Doveton ⁽⁴⁾.
 - RHOMaa – Apparent Grain Density computed from RHOB and NPHI log curves are mapped to the red color range initially from 2.6 to 2.87 gm/cc.
 - Umaa – Apparent Photoelectric computed from PE, RHOB and NPHI log curves are mapped to the green color range initially from 4.3 to 13.6 barnes/cc.
 - GR – Gamma Ray log curve mapped to the blue color range initially from 0 to 150 API.
- The Litho-Density Log Curves without Photoelectric Factor (PE) Colorlith Plot Track, which uses the Bulk Density (RHOB) and Neutron Porosity (NPHI) log curves to compute the RHOMaa curve, the limits suggested by Doveton & Collins ^(2,4).
 - RHOMaa – Apparent Grain Density computed from RHOB and NPHI log curves are mapped to the red color range initially from 2.6 to 2.87 gm/cc.
 - NPHI – Neutron Porosity log curve mapped to the green color range initially from 0.0 to 0.3 PU.
 - GR – Gamma Ray log curve mapped to the blue color range initially from 0 to 150 API.

Color Image Plot Tracks

The Color Image Plot Tracks⁽⁶⁾ uses one curve to plot its magnitude over a 16 color range. The KIMELEON web app sets up 3 color image track types; Porosity, Conductivity and Gamma Ray in both linear and non-linear color scheme. There is only one log curve per type and is automatically selected by a hierarchy depending on the log curves available, which can be modified by the user. The linear color image plot track assigns the color by its magnitude; there are 16 colors within the curve type minimum and maximum value, the minimum and maximum log curve value for each color is divided equally within the overall minimum and maximum values, i.e. for gamma ray the overall minimum and maximum value is 0 and 150 API respectively, divide the range by 16 and you get about 9.375 so the 1st color interval is from 0 to 9.375, the 2nd color interval is from 9.375 to 18.75 etc. up to 16th color interval set from 140.625 to 150. The non-linear color image plot track first sorts all the data for the selected log curve by magnitude dividing all the log values equally among the 16 color intervals and then resorts the log curve values by depth. This value is only dependent on the depth range and not on the curve limits. The KIMELEON web app has set up three color image plot types as follows,

1. Porosity Log Color Image Plot Track, the user can select from the following curve types,
 - a. DPHI – Density Porosity
 - b. NPHI – Neutron Porosity
 - c. SPHI – Sonic Porosity
2. Conductivity (1000.0/Resistivity) Log Color Image Plot Track, the user can select from the following log curve types,
 - a. RES – General Resistivity
 - b. RDEP – Deep Resistivity
 - c. RMED – Medium Resistivity

- d. RSHAL – Shallow Resistivity
 - e. ILD – Deep Induction Resistivity
 - f. ILM – Medium Induction Resistivity
 - g. SFLU – Spherically Focused Resistivity
 - h. LL – Deep Laterolog Resistivity
 - i. MLL – Micro Laterolog Resistivity
 - j. LL8 – Shallow Laterolog Resistivity
 - k. LN – Long Normal Resistivity
 - l. SN – Shallow Normal Resistivity
 - m. MNOR – Micro Normal Resistivity
 - n. MSFL – Micro Spherically Focused Resistivity
 - o. MINV – Micro Inverse Resistivity
 - p. AHT10 – Array Induction Resistivity -10
 - q. AHT20 – Array Induction Resistivity -20
 - r. AHT30 – Array Induction Resistivity -30
 - s. AHT60 – Array Induction Resistivity -60
 - t. AHT90 – Array Induction Resistivity -90
3. Gamma Ray Log Color Image Plot Track, the user can select from the following log curve types,
- a. GR – Gamma Ray
 - b. CGR – Gamma Ray Minus Uranium

References:

- (1) Color display of well logs, Peter L. Briggs, Mathematical Geology, Volume 17, Number 4, May 1985
- (2) Color Images of Kansas Subsurface Geology from Well Logs, D. R. Collins and J. H. Doveton, Computer & Geosciences, Vol. 12, No. 4B, pp.519-526 1986
- (3) Visualization of Subsurface Geology from Wireline Logs, David R. Collins, Digital Mapping Techniques '98-Workshop Proceedings U. S. Geological Survey Open-File Report 98-487
- (4) The Dakota Aquifer Program Annual Report, FY89 Kansas Geological Survey, Open-File Rept. 90-27 Annual Report, FY89-Appendix B
(http://www.kgs.ku.edu/Dakota/vol3/fy89/app_b.htm)
- (5) Composite Colour Display of Spectral Gamma-Ray Logs, by Matt Hall, Canadian Well Logging Society, Dec 2005, v.24
- (6) Geological Well Logs Their use in Reservoir Modeling, by Stefan M. Luthi, pg 84-86, ISBN 3-540-67840-9, © 2001 Springer-Verlag Berlin

Changing the Colorlith Plot Track Limits & Log Curve Limits



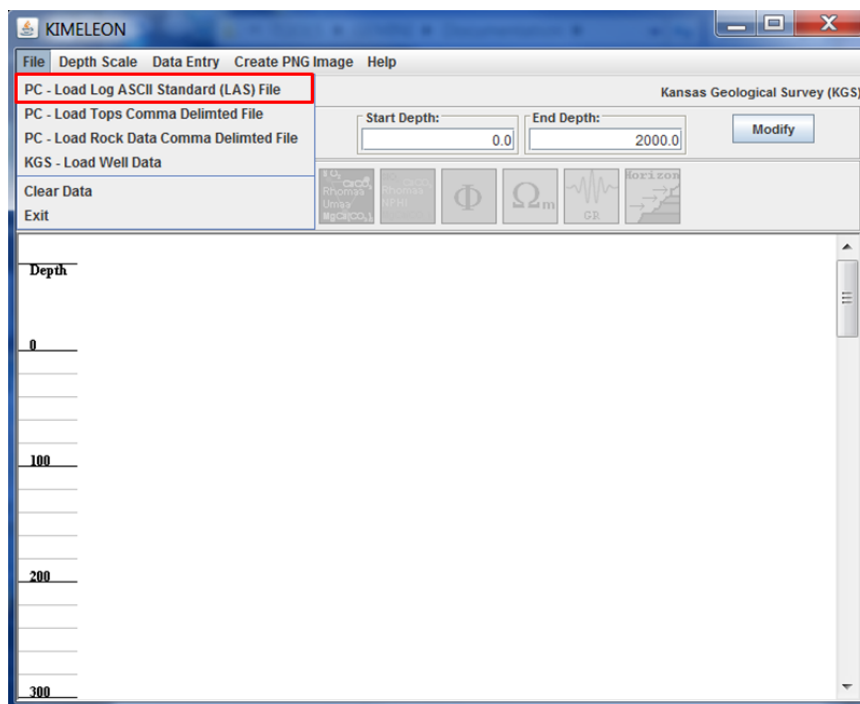
Each of the image icons above opens a specific log type plot panel that allows the user to change the limits of the log curve plots as well as the limits controlling the colorlith and color image plot tracks. This example will use the Braun 1 Log ASCII Standard (LAS) version 3.0 file, which can be downloaded from the following URL's,

Well Data: Braun 1, Ellis County, Kansas

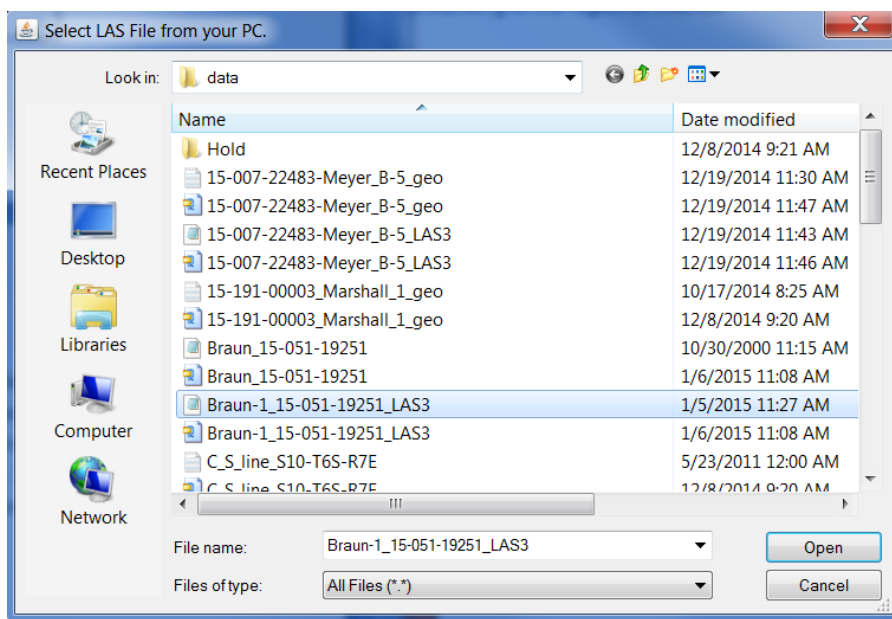
Type	ASCII Text Files
LAS 3.0	http://www.kgs.ku.edu/Gemini/Tools/documentation/Braun-1_15-051-19251_LAS3.las

Type	Zip Files
LAS 3.0	http://www.kgs.ku.edu/Gemini/Tools/documentation/Braun-1_15-051-19251_LAS3.zip

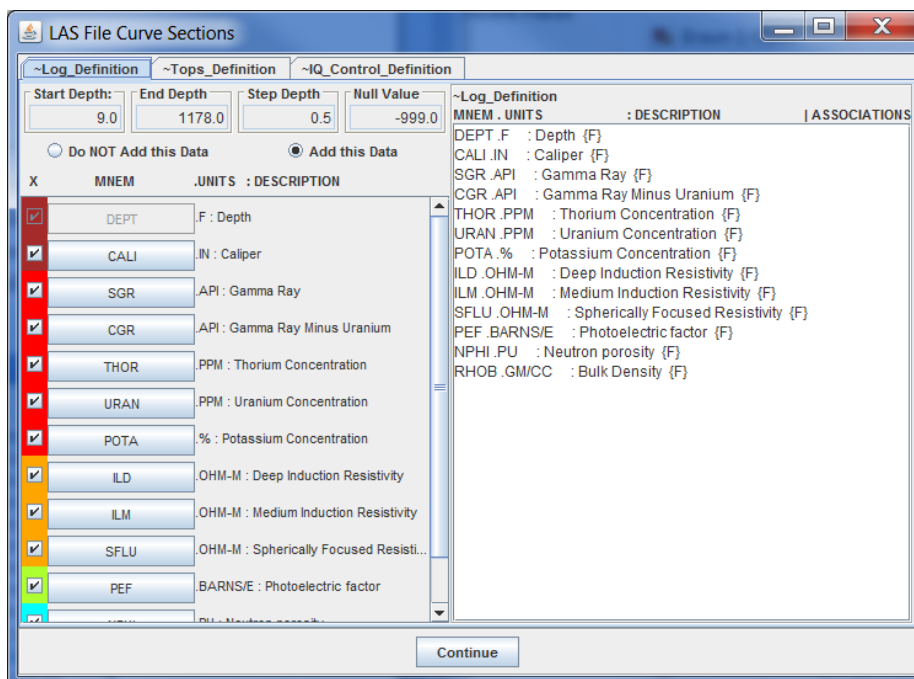
Download either the ASCII Text Files directly or the Zip files extracting the contents into a directory. The problem with the ASCII Text Files being downloaded directly from a web page is that the web page will alter the contents so it does not retain the basic structure and add HTML text to the file. The preferred method if you have Zip or WinZip is to download the zip files to your PC and extract. The Braun-1_15-051-19251_LAS3.las file contains both the Log & Tops Data.



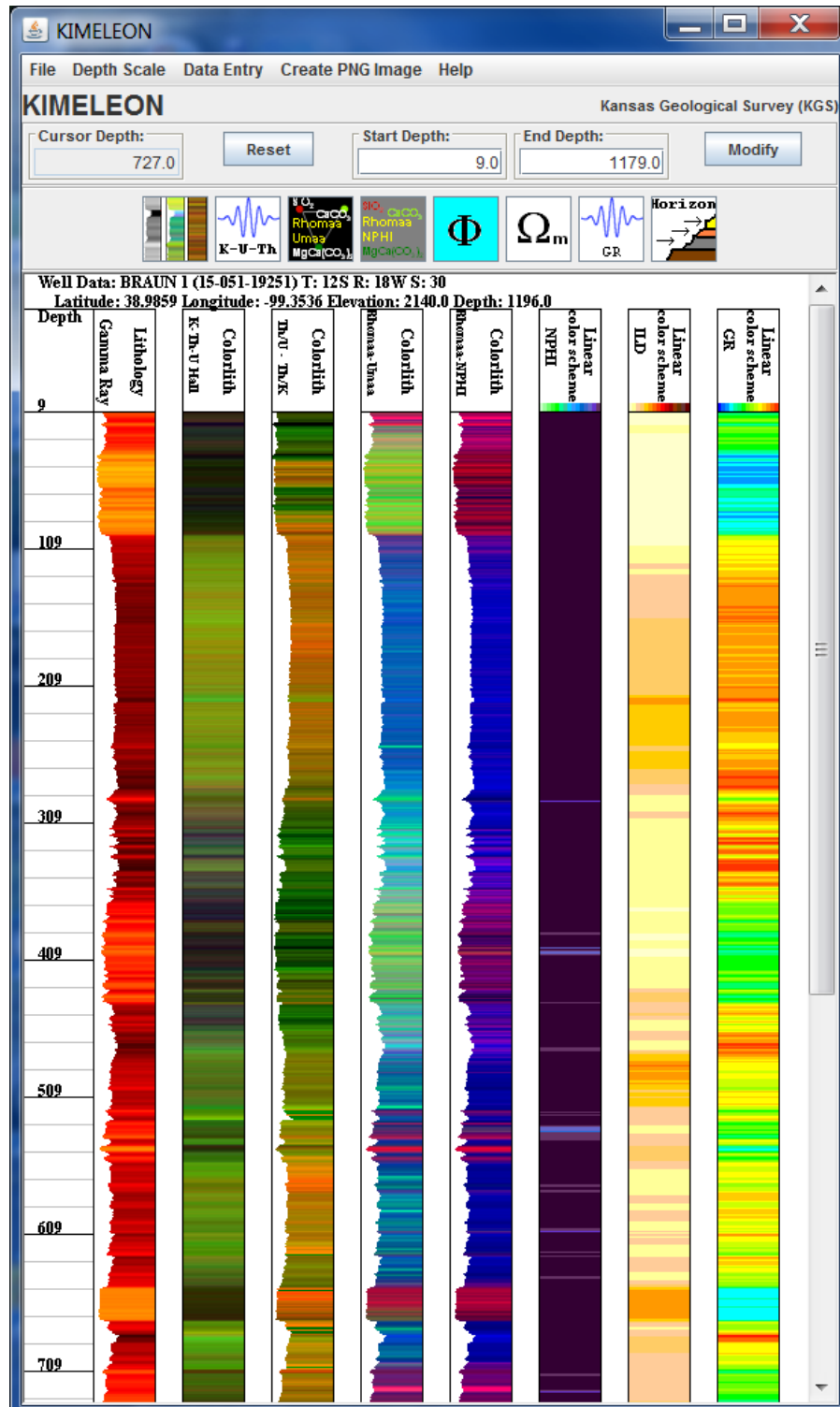
Click on the “File” Menu in the menu bar at the top of the “KIMELEON” Dialog, which will display a list of menu options to import well data into the web app. Select the “PC - Load ASCII Standard (LAS) File” menu option. This will display the “Select LAS File from your PC” Dialog. This dialog allows the user to search their PC for the file of interest. In this example it is the LAS version 3.0 file Braun-1_15-051-19251_LAS3.las, highlighted below. Select the Open button to display the “LAS File Curve Sections” Dialog.



The “LAS File Curve Sections” Dialog allows the user to map unknown LAS Curve Mnemonics to the KGS “Standard” Curve Mnemonics so they will be plotted in the LAS File Viewer Plot.



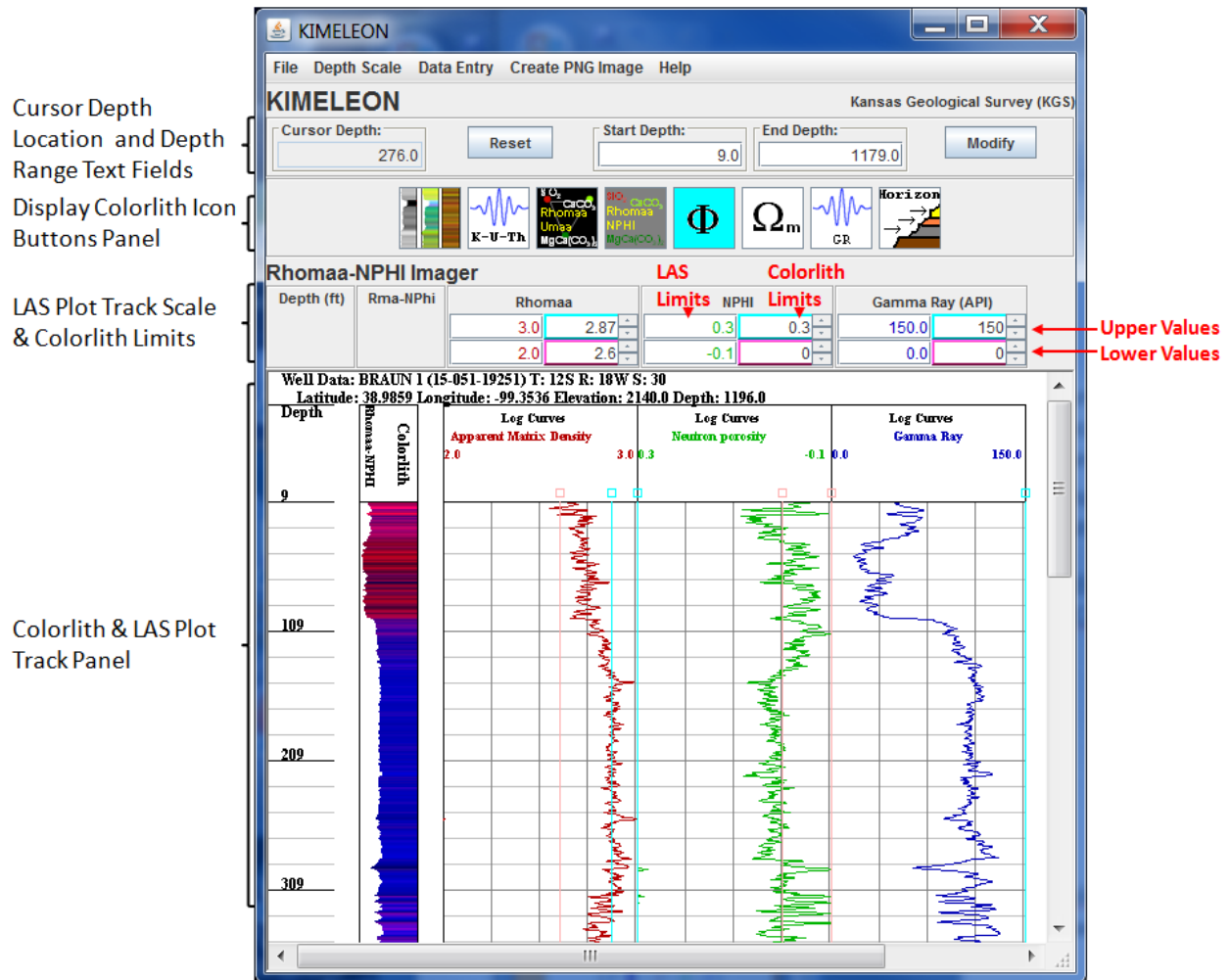
This program reads the “LAS Tool Curve Mnemonics map to KGS Standard Mnemonics” XML File (http://www.kgs.ku.edu/software/gemini/data/las_standard_tools.xml), which will automatically maps the Curve Mnemonics from the LAS file to one of 31 KGS “Standard” Curve Mnemonics. This file has all the curve correctly mapped so select the “Continue” button to load the log and tops data into the KIMELEON web app.



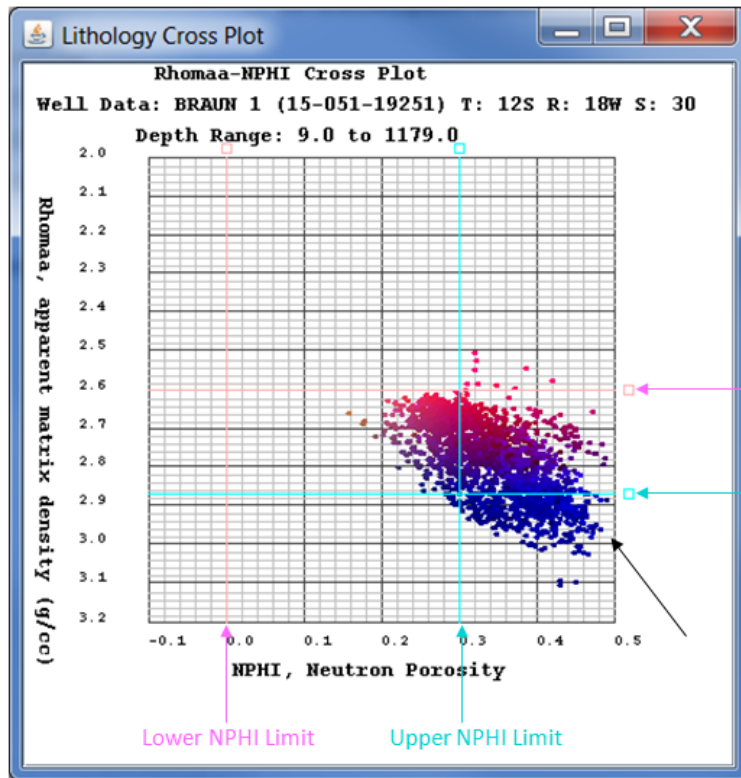


Rhoma-NPHI Icon Button

Select the “Rhoma-NPHI” Icon Button to display the “Rhoma-NPHI Image” Panel and the Rhoma-NPHI 2D Plot Panel.



The 2D Plot Panel only displays for the “Rhoma-Umaa” and “Rhoma-NPHI” Icon Image buttons. The Colorlith limits colors cyan and pink match the Colorlith Limits text fields.



Colorlith Curve Limits

- Cyan color is the upper limit.
- Pink color is the lower limit.

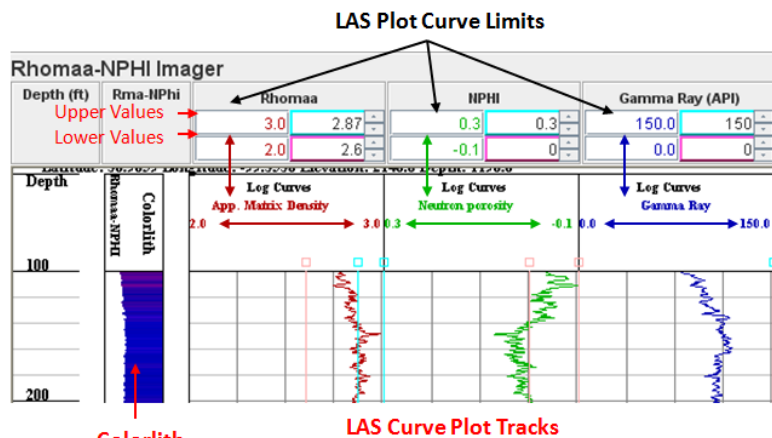
If curve data falls outside of curve limit range the curve data is redefined to be the maximum or minimum of the limit.

Color of data is the (Red, Green, Blue) equivalent of the data (Rhomaa, NPHI, Gamma Ray) values respectively within the limits of the colorlith color values.

The points displayed in the plot fall within the depth range specified on the Kameleon Control Panel.

LAS Plot Curve Limits

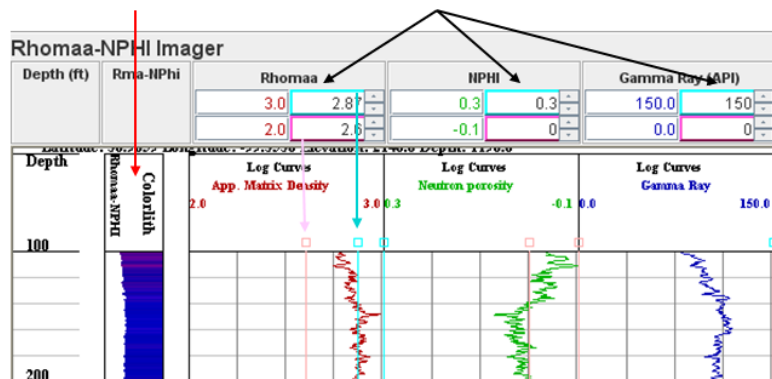
- Each Log Curve has a default curve limit, user can modify the limits.
- Color of the Plot limit matches the color of the Log Plot & Labels.
- Color of the Log Curve represents the color used in the tri-color RGB colorlith presentation, i.e.
 - Red – Rhomaa Log Curve
 - Green – Neutron Porosity Log Curve
 - Blue – Gamma Ray Log Curve



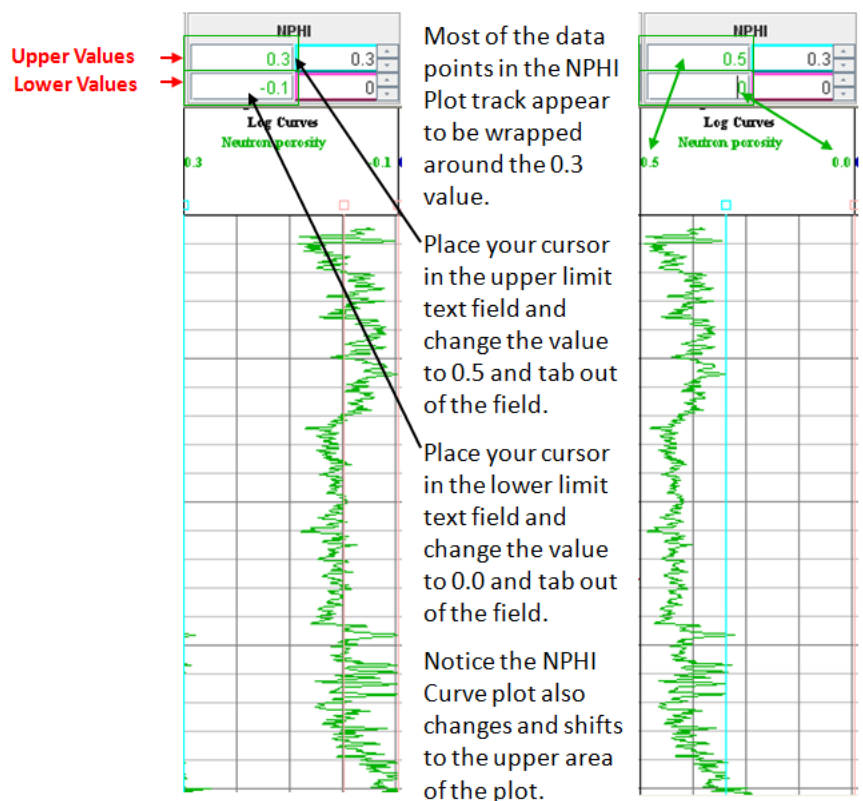
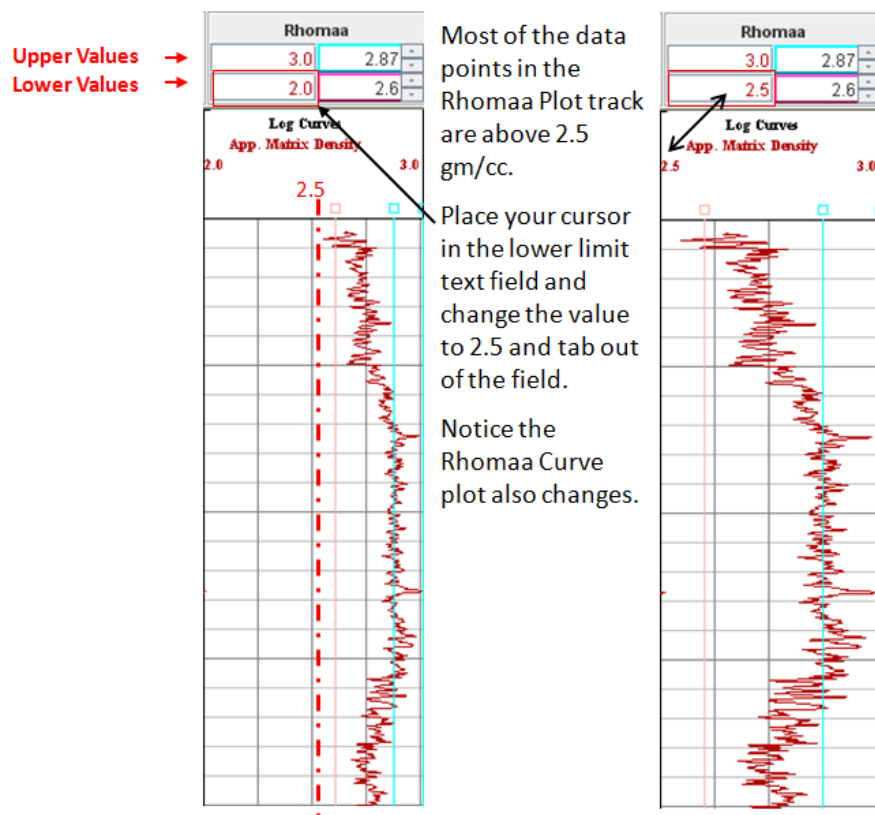
Colorlith Curve Limits

Colorlith Curve Limits

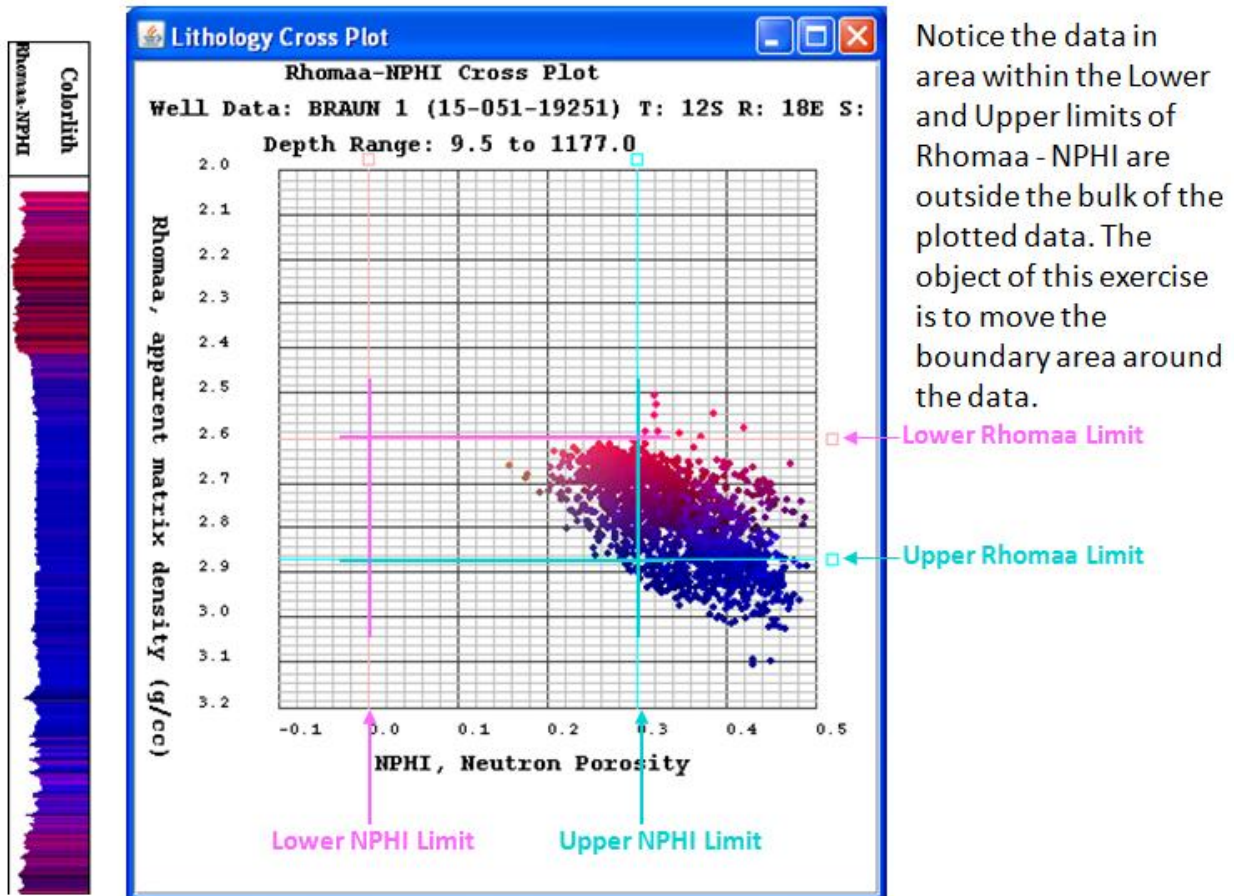
- Each Log Curve has a default colorlith plot limit, user can modify the limits.
- Border Color of the colorlith text field matches the square and line color on the LAS Plot Track.
- Cyan color is the upper limit.
- Pink color is the lower limit.

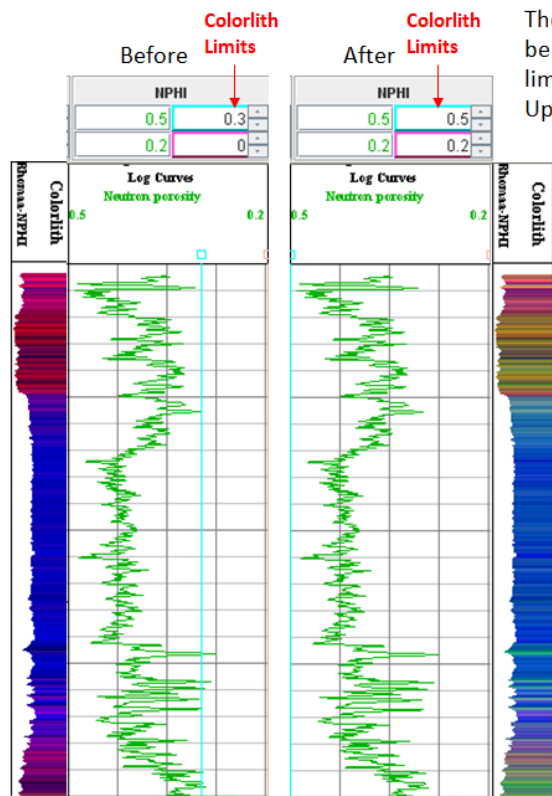


Changing LAS Plot Limits

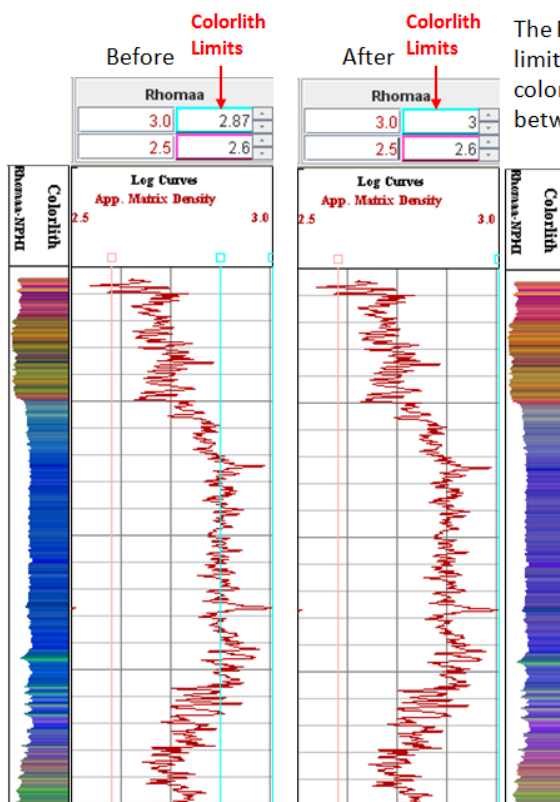
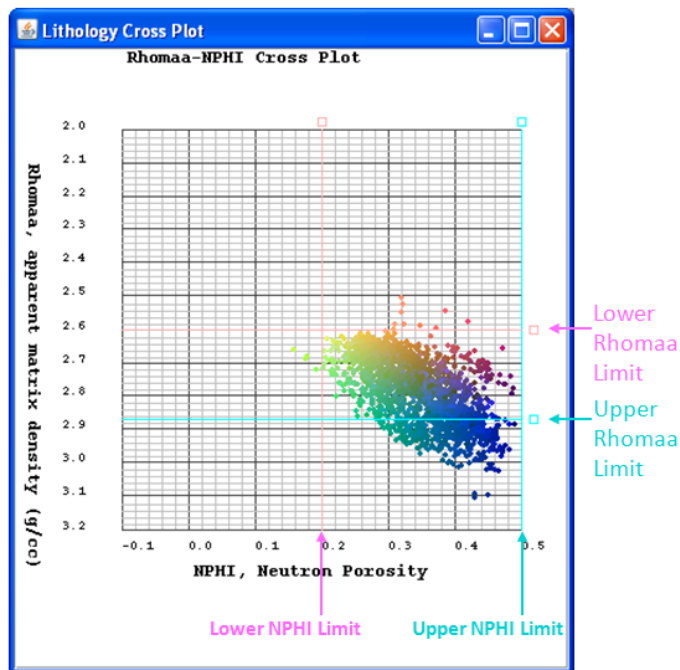


Changing Colorlith Limits

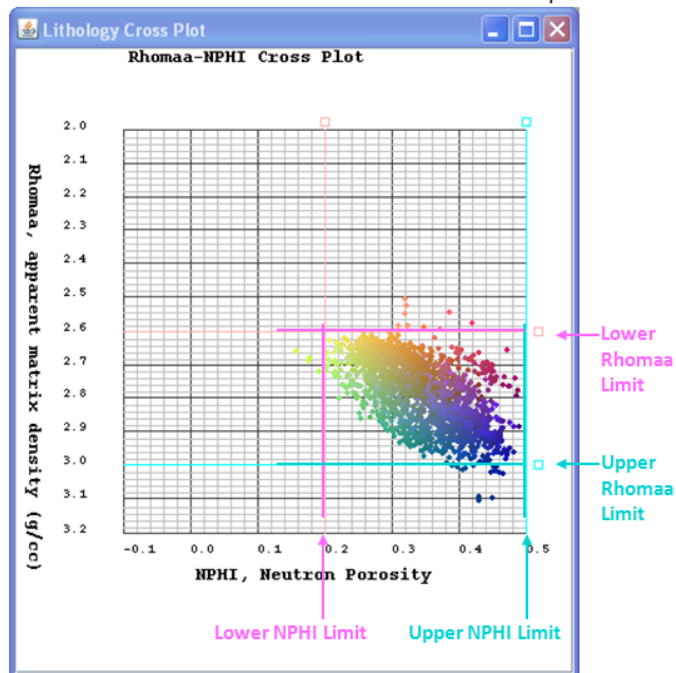




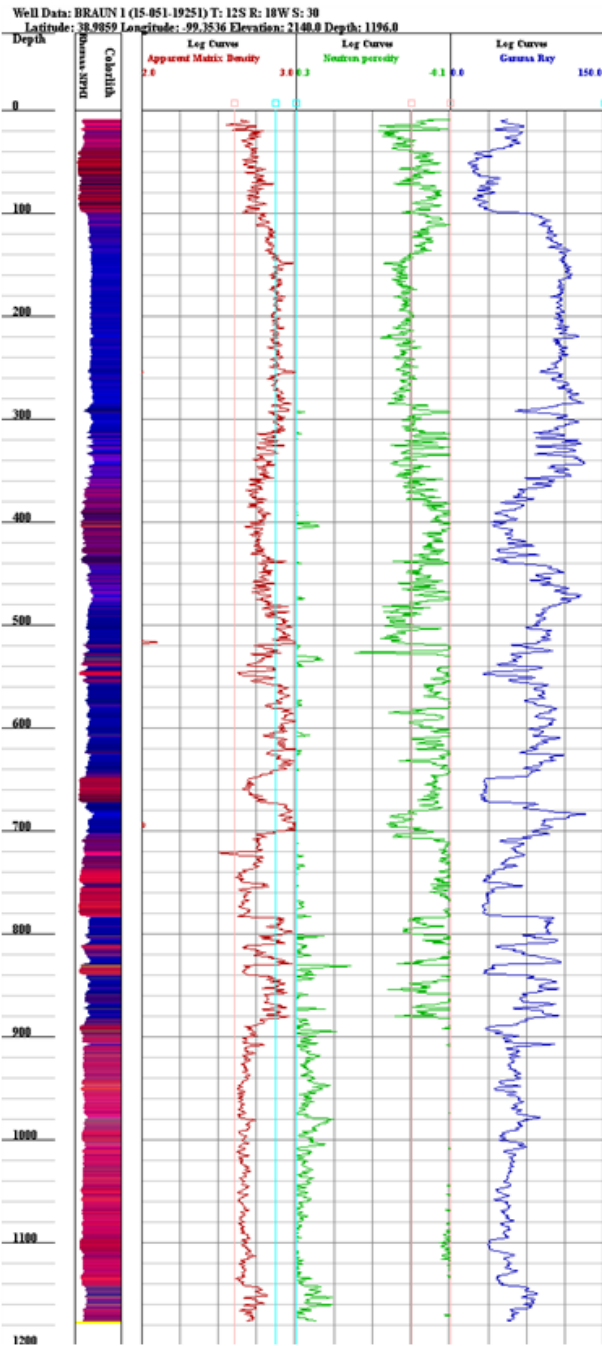
The Rhomaa-NPHI Cross Plot illustrates the cluster of data points to be above 0.2 PU and below 0.5 PU. Change the limits in the colorlith limits text fields and notice in the Cross Plot that the Lower and Upper NPHI Limits now encloses the data points in NPHI space.



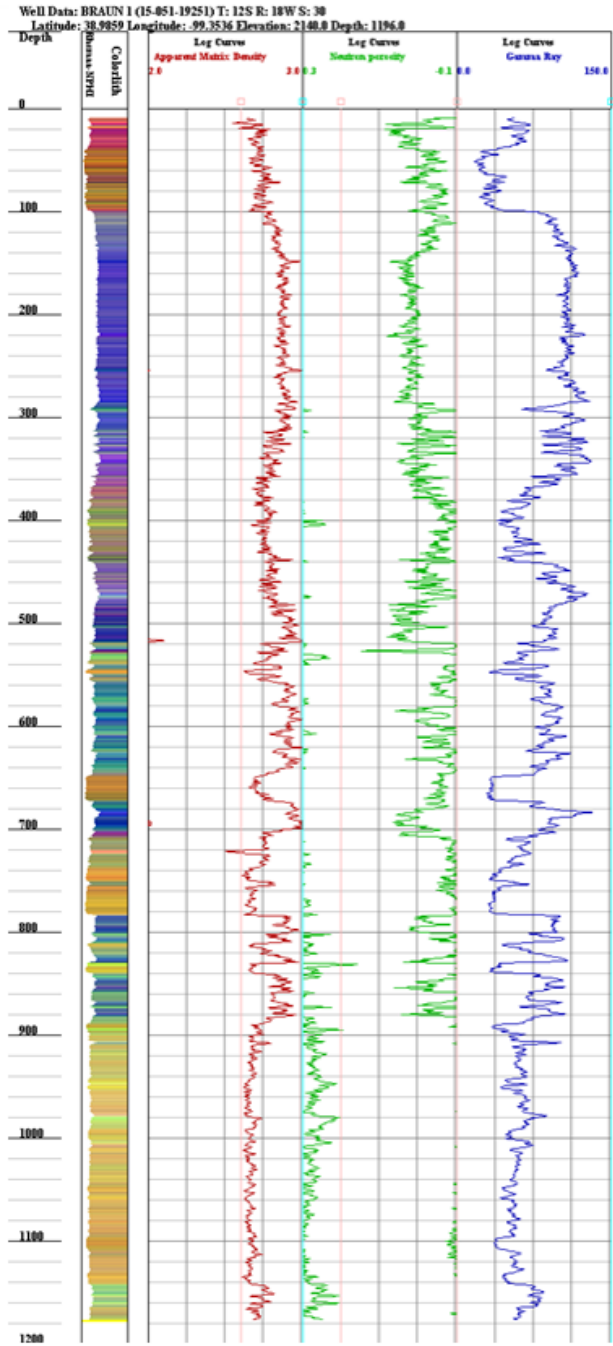
The Rhomaa-NPHI Cross Plot illustrates that the Rhomaa curve limits are only about half of the data points. Increase the upper colorlith curve limit to 3.0 and notice that the boundary area between NPHI and Rhomaa now encloses most of the data points.



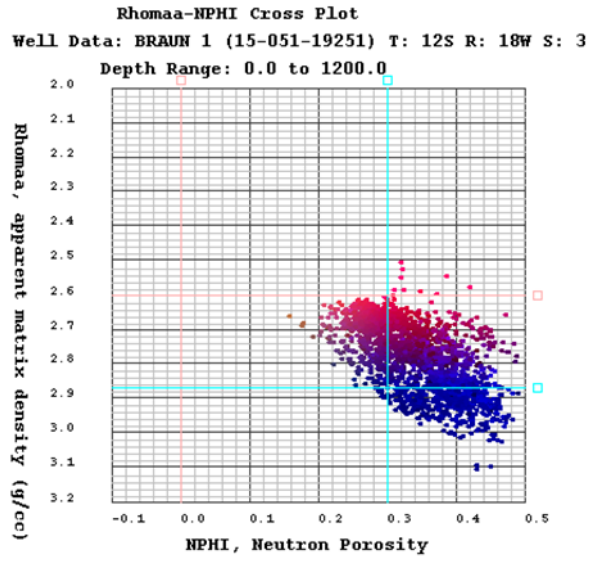
Before Colorlith Limit Changes



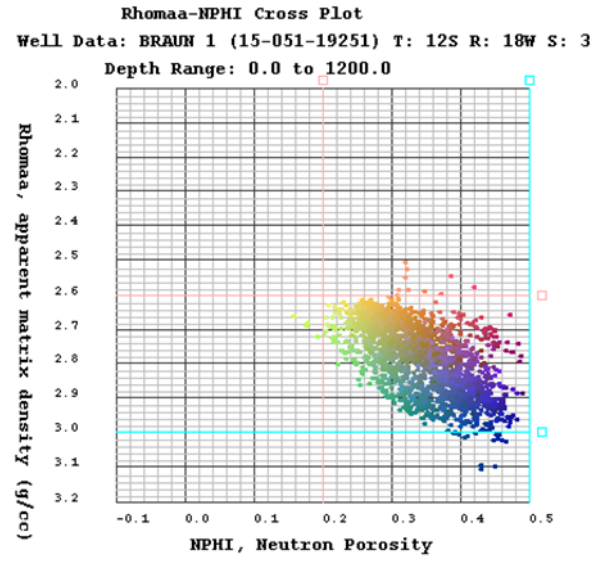
After Colorlith Limit Changes



Before Colorlith Limit Changes



After Colorlith Limit Changes



Change Depth Range & Scale

This program allows the user to change the viewing depth range for the plot and to change the feet / inch plot scale to expand or contract the Kimeleon plot.

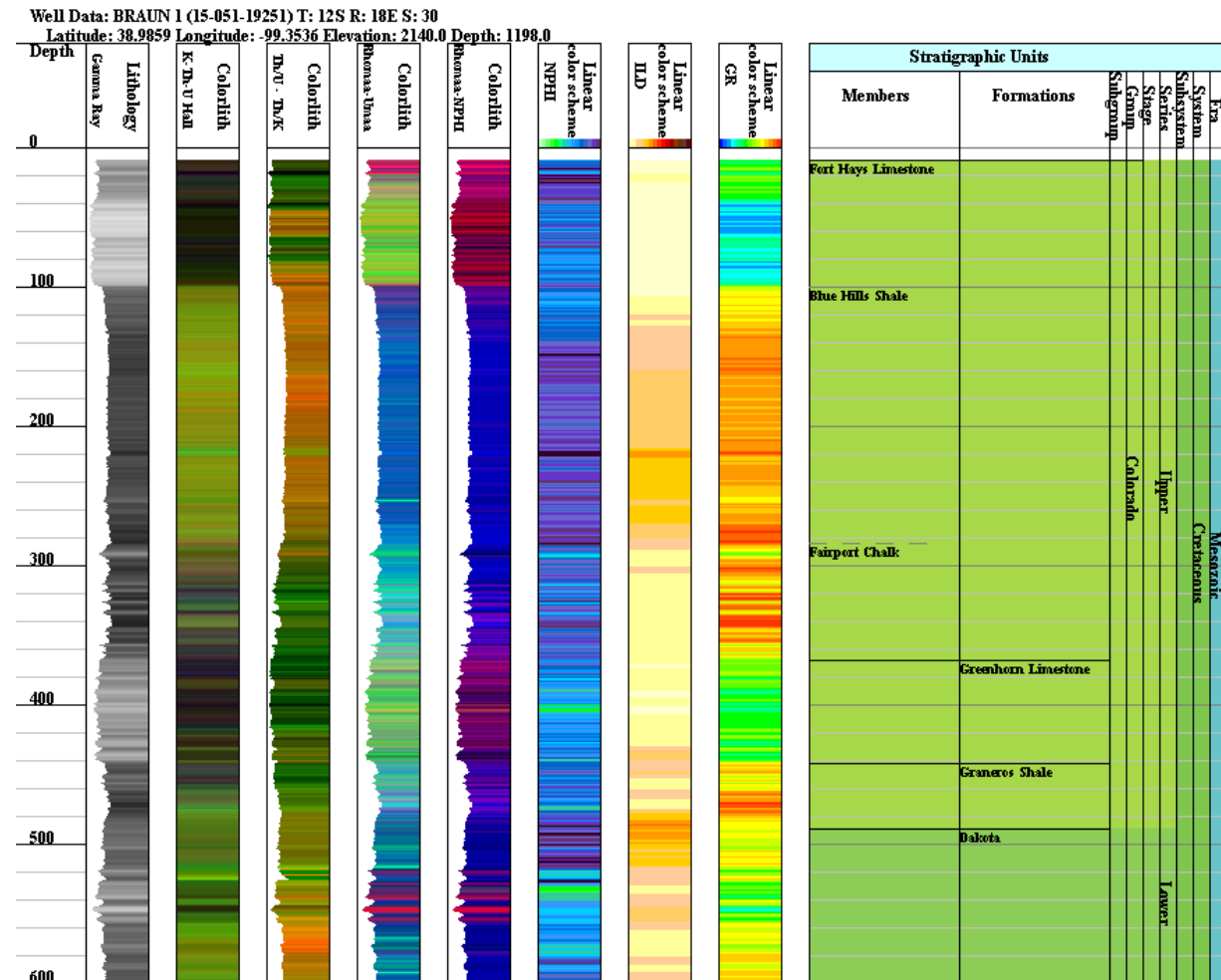
Go to the top of the Kimeleon Frame. This example the Kimeleon Plot of the Braun 1 will be changed from 9.5 Start Depth, 1177.0 End Depth to 0.0 Start Depth, 600.0 End Depth.
Select the "Modify Depth" button.

Before

Reset	Start Depth: <input type="text" value="9.5"/>	End Depth: <input type="text" value="1177.0"/>	Modify
-------	--	---	--------

After

Reset	Start Depth: <input type="text" value="0.0"/>	End Depth: <input type="text" value="600.0"/>	Modify
-------	--	--	--------



Before

Depth Scale

- ☐ 2 ft / in
- ☐ 5 ft / in
- ☐ 10 ft / in
- ☐ 20 ft / in
- ☐ 50 ft / in
- ☒ 100 ft / in
- ☐ 200 ft / in

This example the Kimeleon Plot of the Braun 1 will change the depth scale of 100 ft / in to 200 ft / in. Notice that the "Depth Scale & Range" Panel shows the Depth Scale to be 100 ft / in.

Go to the "Depth Scale" Menu on the "Kimeleon Control" Frame. Select the "200 ft / in" radio button.

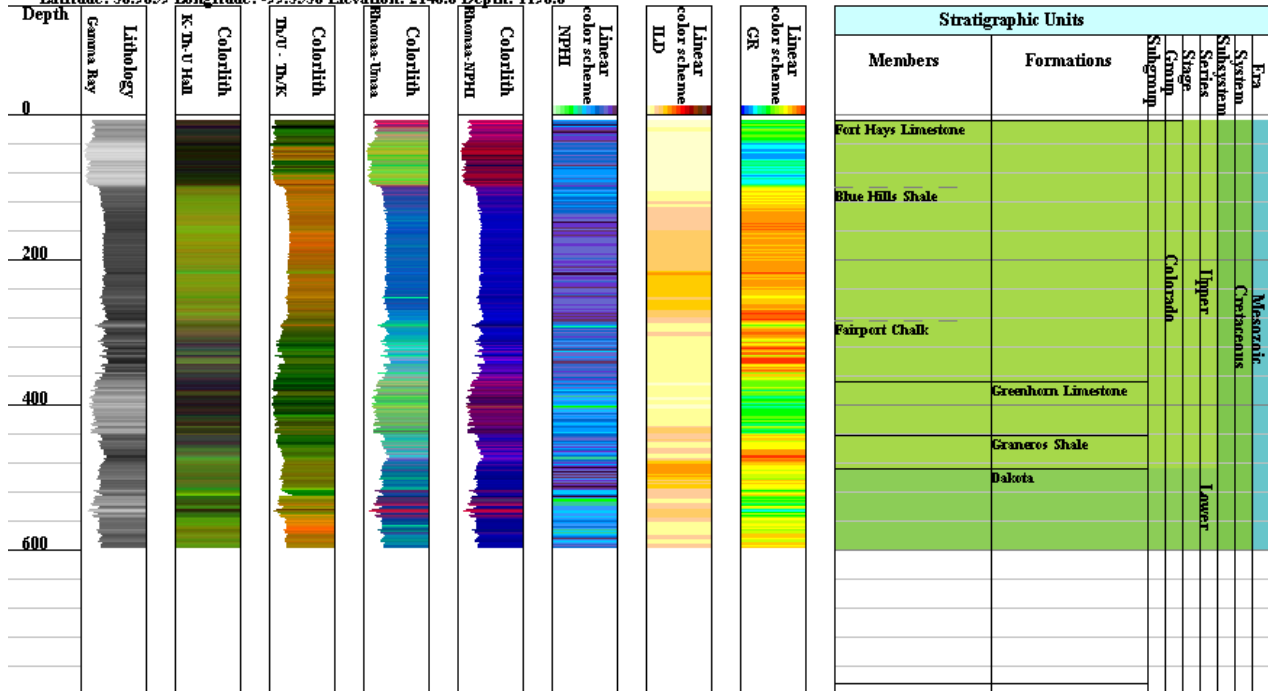
Notice that the "Depth Scale & Range" Panel shows the Depth Scale to be 200 ft / in and the Kimeleon plot has contracted.

After

Depth Scale

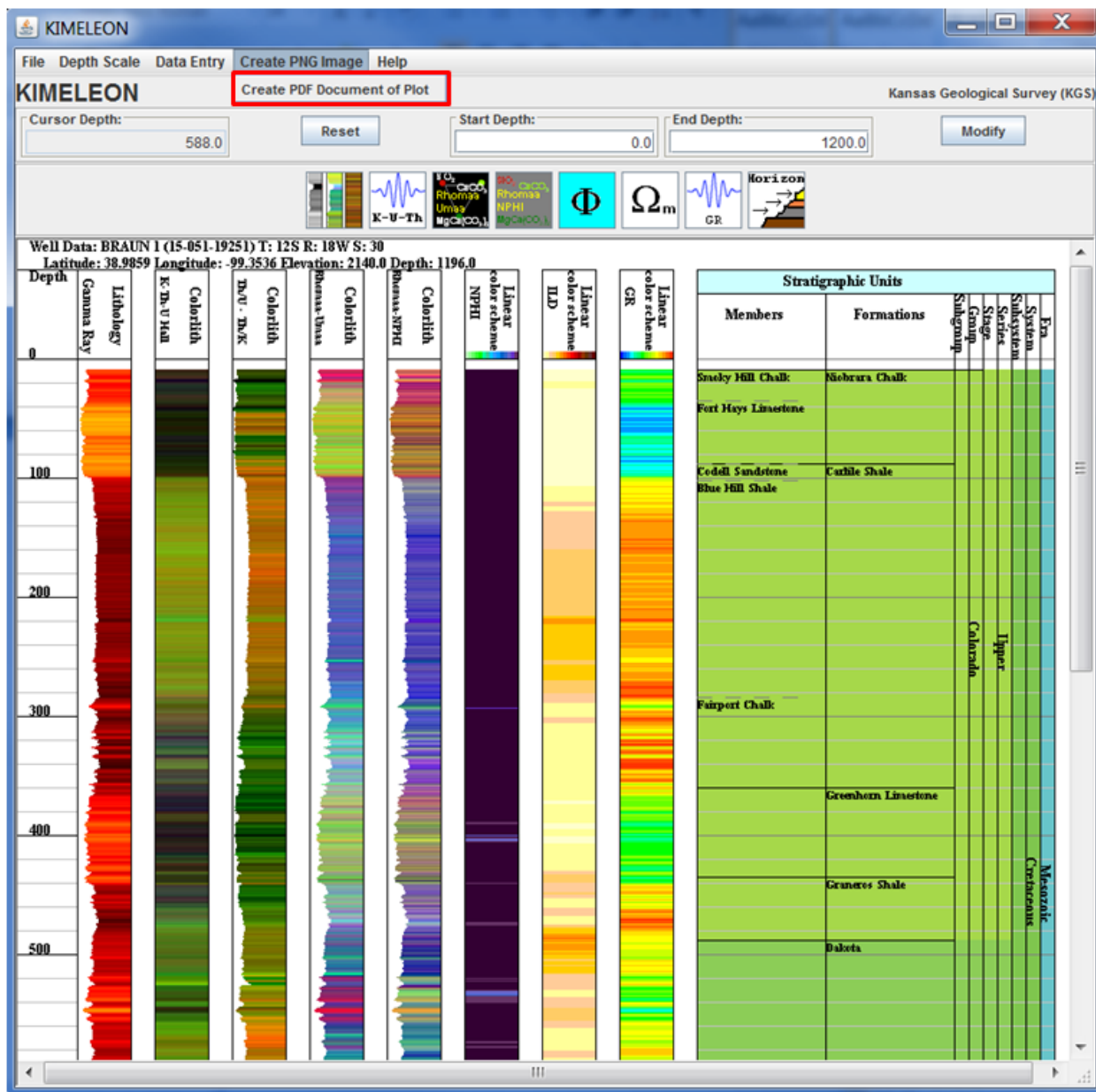
- ☐ 2 ft / in
- ☐ 5 ft / in
- ☐ 10 ft / in
- ☐ 20 ft / in
- ☐ 50 ft / in
- ☐ 100 ft / in
- ☒ 200 ft / in

Well Data: BRAUN 1 (15-051-19251) T: 12S R: 18E S: 30
 Latitude: 38.9859 Longitude: -99.3536 Elevation: 2140.0 Depth: 1198.0

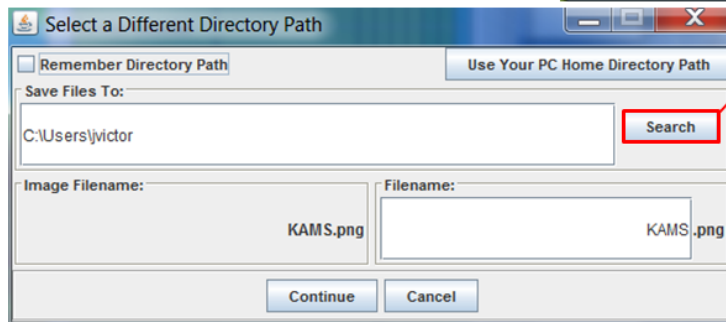


Creating a Portable Document Format (PDF) Document of Plots

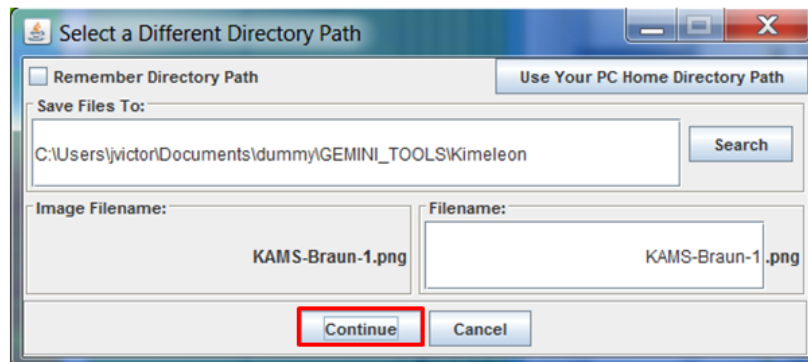
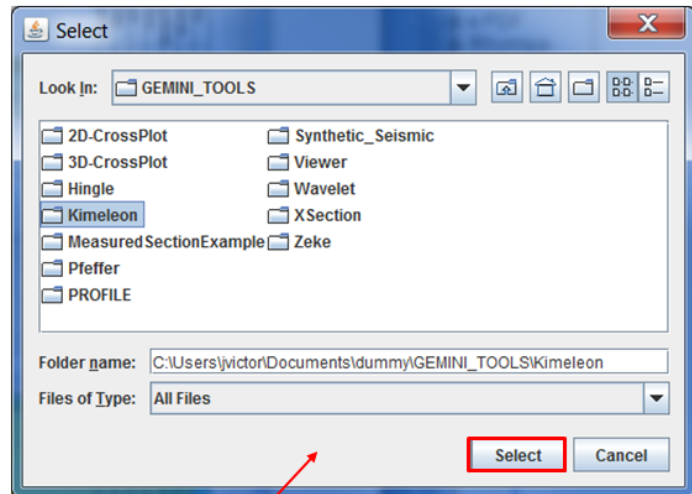
Select the “Create PNG Image” Menu and then select the “Create PDF Document of Plot” menu option, which will display the “Select a Different Directory Path” dialog which allows the user to place the Portable Network Graphics (PNG) image file anywhere on their PC. Also when the user selects the “Create PDF Document of Plot” menu option the panel that is displayed is the PNG image that is being created. Only the “Rhomaa-Umaa” & “Rhomaa-NPHI” icon buttons will also create a 2D Plot PNG image with the Plot Panel.



Click on the “Search” button to display the “Select” dialog that allows the user to search the PC for the directory path to save the PNG Files.

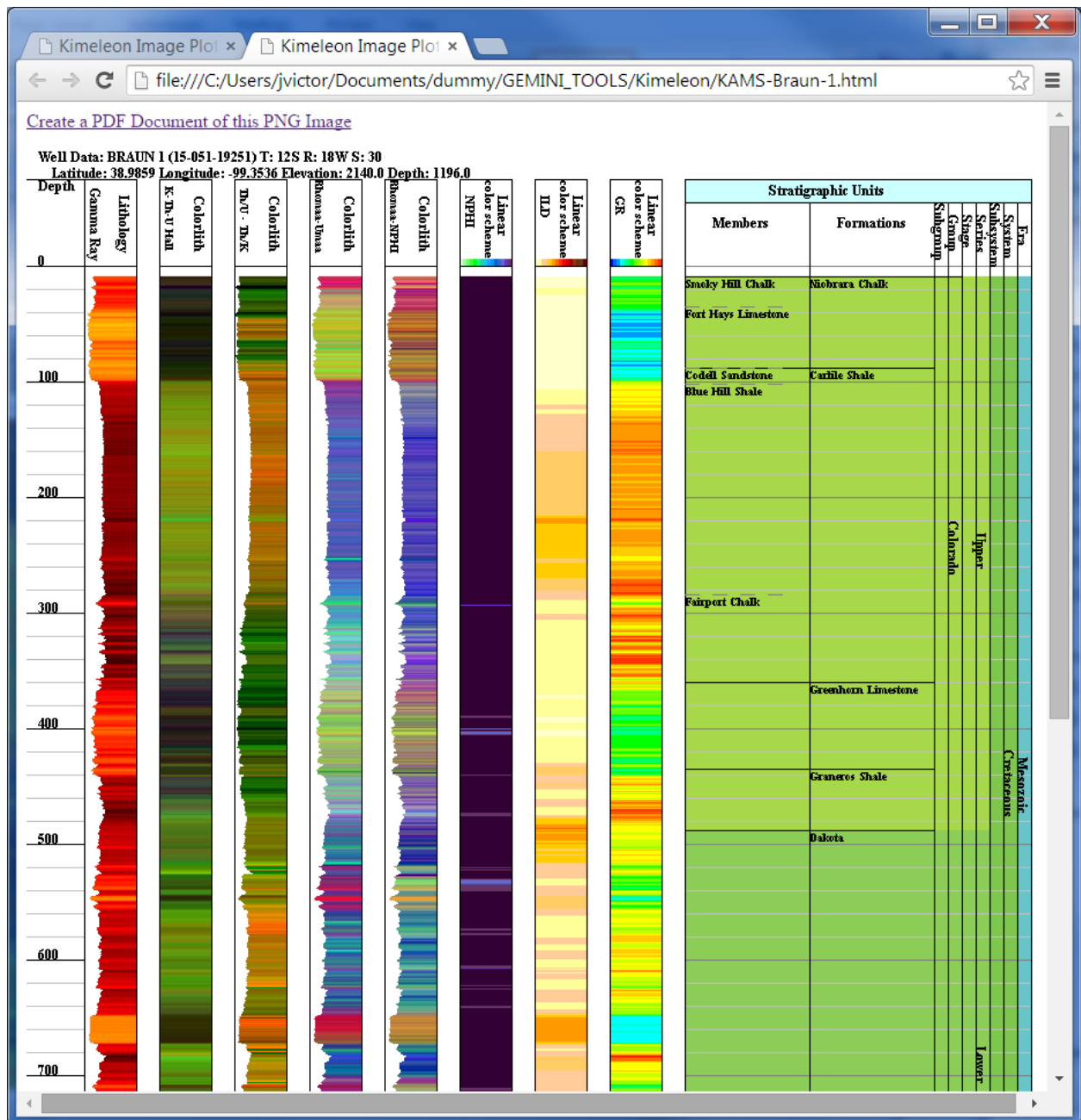


Click on the “Select” button to transfer the directory path to the “Save Files To” text field on the “Select a Different Directory Path” dialog.



The KAMS Prefix identifies the Mnemonic of the panel that is being saved as a PNG file. “-Braun-1” is being added to “KAMS” to identify the Well with the image type.


Select the “Continue” button to create the PNG file.



When the PNG image is created a HTML document is also created to show the created PNG image with a link at the top that allows the user to launch the PDF (Portable Document Format) Applet . Click on the “Create a PDF Document of this PNG Image” URL link.

Kimeleon Image Plot x Create PDF Document x

www.kgs.ku.edu/PRS/Ozark/TYPE_LOG/PDF.html




Create PDF (Portable Document Format) from Image Applet

This applet allows the user to create a PDF (Portable Document Format) from a PNG (Portable Network Graphics), JPEG (Joint Photographic Experts Group) or GIF (Graphics Interchange Format) image file.

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The URL for this page is http://www.kgs.ku.edu/PRS/Ozark/TYPE_LOG/PDF.html

Do you want to run this application?


 **Name:** PDF

Publisher: University of Kansas

Location: <http://www.kgs.ku.edu>

This application will run with unrestricted access which may put your computer and personal information at risk. Run this application only if you trust the location and publisher above.

☐ Do not show this again for apps from the publisher and location above

 More Information Run Cancel

Select the "Run" button to launch the search directory path on your PC dialog to find the PNG that was just created.

