





HyLogger Case Study: Greenbushes (C3DD024)

Pt. 1 Geological Context and Introduction to NVCL Datasets

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National Virtual Core Library (NVCL) Project



Summary

This case study includes a HyLogger data pack for drill hole C3DD024 (Greenbushes Li-Sn-Ta pegmatite deposit, Western Australia) comprising hyperspectral drill core reflectance spectra ("HyLogger data"), geochemistry, and lithology data, as well as a case study report split into several parts. The *.tdg file* allows full access to the data and CSIRO's The Spectral Geologist software (TSG[™]) without a full TSG software license required for this dataset. Part I of the case study report outlines the geological context of the C3DD024 drill hole, providing the necessary background for interrogating the datasets. It will also provide an overview of the datasets available with the case study. Subsequent parts of the case study will aim to interrogate specific geological questions using this dataset, such as white mica and smectite alteration signatures. New applications-based sections will be added periodically to the case study and will use the same datasets.

Geological Context

The Greenbushes Li-Sn-Ta pegmatite deposit is a world class Li mineral asset and occurs ~350 km south of Perth, in Western Australia, in the Southwest Terrane Greenstones of the Yilgarn Craton (Figure 1). The Greenbushes deposit presently supplies more than 30% of the worlds lithium and is the single largest Li hard rock deposit in the world. Development of the Greenbushes Li orebody began in 1983, however, the Greenbushes areas has been mined for tin since 1888 and tantalum since 1984 (Talison Lithium). The resource for Greenbushes was 158.06 Mt of ore at a grade of 2.25 wt% Li₂O for 8.72 Mt of lithium carbonate equivalent in 2015 and has been continuously mined since this time (Partington, 2017).

The Greenbushes pegmatites are rare-element pegmatites belonging to the Lithium-Caesium-Tantalum (LCT) family and occurs in the Archean Western Gneiss Terrane (Figure 1). The pegmatites intruded into the Balingup Metamorphic Belt along the 2.53 Ga Donnybrook-Bridgetown shear zone. This shear zone is NNW trending, 15-20 km wide, and consists of sheared gneisses, amphibolite, and syn-tectonic granitic intrusions (Partington et al, 1995).

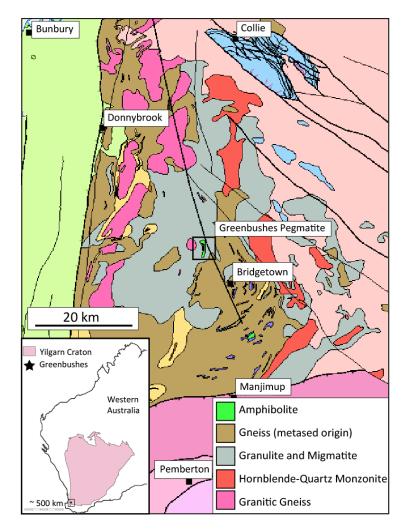


Figure 1 Regional map showing the geology and location of the Greenbushes Pegmatite in the southwest corner of the Yilgarn craton (modified after Duuring, 2020 and GeoVIEW.WA, 2016)

The region has undergone several stages of deformation beginning with the formation of the Bridgetown gneisses at 3.1 Ga, deformation and intrusion of regional granites from 2.61 to 2.58 Ga, followed by intrusion and deformation of the Greenbushes Pegmatite at 2.53 Ga (Partington et al, 1995). The pegmatites were intruded by dolerite dykes at 2.43 Ga resulting in some hydrothermal remobilization followed by several younger deformation and intrusion events. For a detailed summary of the regional deformation history, see Partington et al, 1990, 1995 and Partington, 2017. The mineralization at the Greenbushes Pegmatite is unusual in that it occurs in a high-temperature and high-pressure metamorphic terrane but was intruded 50 Ma after regional granite intrusion and so is not associated with any known proximal fractionated granites (Partington, 2017).

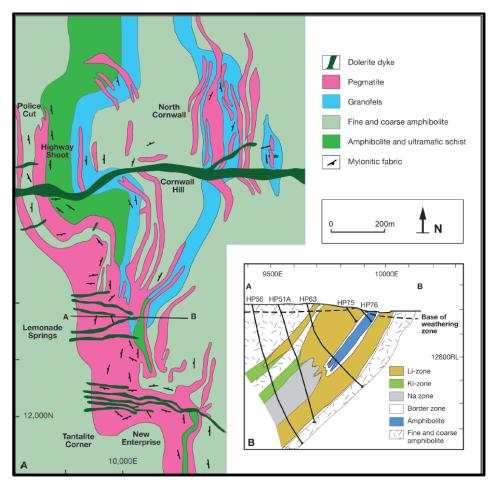


Figure 2 Mine scale geology of the Greenbushes with an inset cross section (A-B) showing the pegmatite zonation and distribution of Lithium in the main ore zone (From Partington, 2017)

The pegmatites occur within shear zones at the contact between major sequences of country rock (granofels, ultramafic schist, amphibolite) and are crosscut by dolerite dykes. The pegmatites are mineralogically zoned and consist of a series of linear intrusions which are 2-3 km in strike and 10 to 300 m thick (Figure 2). The main Li zone is over 2 km long and the main ore mineral is spodumene, a Li-pyroxene (LiAl(SiO₃)₂), which is mined in open pits from fresh, unweathered zones of pegmatite and often accounts for up to 50% of the rock volume (Partington, 2017). The mineral zonation observed in the Greenbushes pegmatites is unusual with the most Li-rich zones occurring in the hanging wall and footwall zones, surrounding the albite zone. In general, the albite zone hosts the highest-grade tin and tantalum ore shoots (Figure 2, Partington, 2017).

Drill Hole Context

Drill hole C3DD024 was drilled from within the Central Lode open pit and collared in fresh pegmatite (Table 1). It encounters over 200 m of pegmatite in the main lithium ore zone and was drilled to better constrain boundaries of the pegmatite resource (Hancock and Wawryk, 2020). The drill hole also intersects minor (<1 m wide) dolerite units and finishes in ~10 m of amphibolite (Table 2; Figure 3; Figure 4).

Table 1 Drill hole details (Hancock and Wawryk, 2020)

Hole	Lat (°E)	Long (°s)	Elevation	Inclination (°) ⁵	Azimuth (TN) ⁶	EOH
C3DD024	116.0651289	-33.85920461	205.186	-59.7	077.2	246.5 m

Table 2 Summary of the geology log for C3DD024 (Hancock and Wawryk, 2020)

From	То	Lithology
0	153.63	Pegmatite
153.63	154.41	Dolerite
154.41	156.07	Pegmatite
156.07	159.92	Dolerite
159.92	182.63	Pegmatite
182.63	182.82	Dolerite
182.82	196.19	Pegmatite
196.19	197.49	Dolerite
197.49	201.28	Pegmatite
201.28	201.61	Dolerite
201.61	218.79	Pegmatite
218.79	218.91	Dolerite
218.81	237.77	Pegmatite
237.77	246.5	Amphibolite

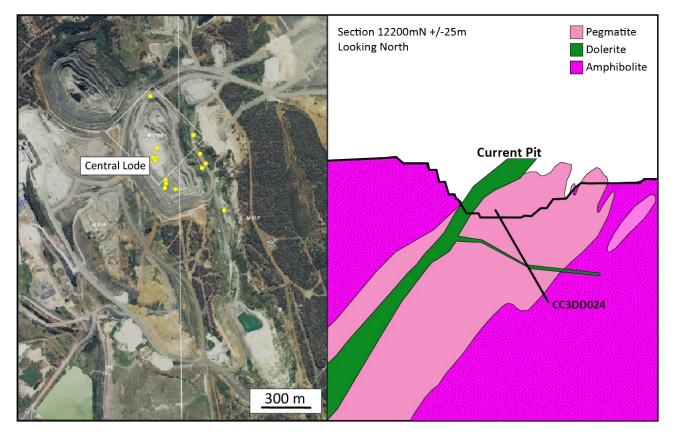


Figure 3 Arial photo of the Greenbushes mine with 2014 drill collars marked as yellow circles and cross section of the Central Lode showing the position of drill hole CC3DD024 (modified from GSWA WAMEX Report A104943 and Talisan Lithium Pty Ltd)

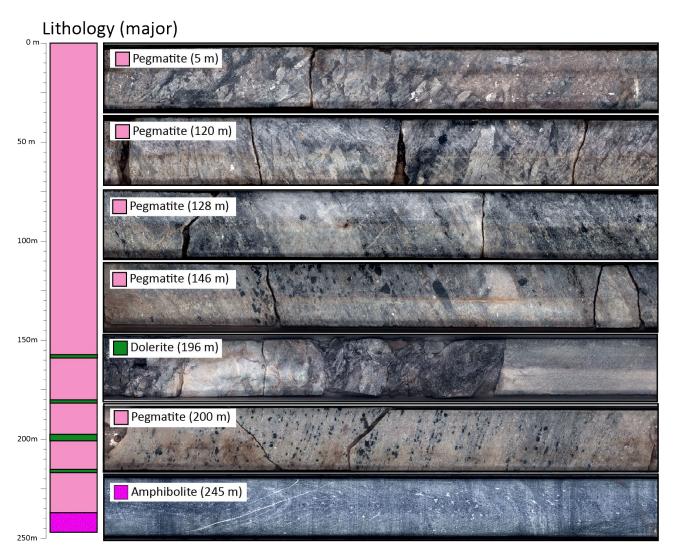


Figure 4 Summary of the lithologies intersected in drill hole C3DD024 with HyLogger photos showing the variability in the pegmatite down hole from coarse grained spodumene at the top with to less spodumene rich to more feldspar rich with notable tourmaline towards the bottom.

Introduction to the Case Study Datasets

The TSGTM file in this case study dataset is provided as a *.tdg file* and includes HyLogger data (i.e. reflectance spectra, high-resolution RGB imagery, laser line profilometry) as well as some imported geochemistry and lithology information (visual logging). To view the HyLogger data, open the file ending in *.tdg* in TSGTM.

This TSG[™] file was generated by the HyLogger-3 system during analysis of the core at the Geological Survey of Western Australia (GSWA) using the method described by Hancock and Huntington (2010), Hancock et al. (2013) and Schodlok et al. (2016). Reflectance spectra were collected over the visible-to-near-infrared (VNIR 380– 1000 nm), short-wave infrared (SWIR 1000– 2500 nm) and thermal infrared (TIR 6000– 14 500 nm) wavelength ranges and high-resolution (0.1 mm pixel) digital colour photographs of core were obtained concurrently using a built-in line-scan camera. The reflectance spectra were automatically resampled to 8 nm spectral sampling and 8 mm spatial resolution using TSG[™].

The results discussed Part I of this case study are primarily based on The Spectral Assistant algorithm (TSA; Berman et al., 2017) outputs which are derived in the TSG[™] software (e.g. Figure 4). TSA is an algorithm for automated spectral unmixing which uses a training spectral library to match the spectrum against a single mineral or model a simulated mixture of 2-4 minerals that most closely resembles that of the input spectrum (Berman et al., 2017). TSA mineralogy outputs are one of the most common outputs derived from hyperspectral data using TSG[™] and for this dataset has been derived using expert input and domain knowledge. The SWIR and TIR spectral ranges are sensitive and diagnostic for different mineral groups, and so the SWIR and TIR TSA result are complementary and should be considered together (Figure 5).

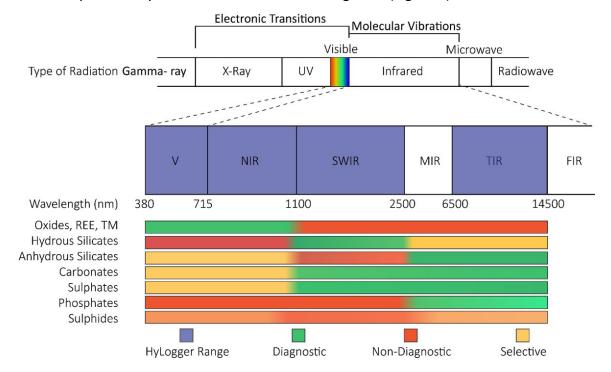


Figure 5 Summary of the Hylogger spectra range and the mineral groups for which the VNIR, SWIR, and TIR are diagnostic, non-diagnostic, and selective.

There are also several scalars discussed in the later parts of this case study (e.g, 2200D). Scalar is the term used by TSG[™] to refer to any set of calculated values related to loaded spectral data. The scalars applied are pre-written, well-established, and in most cases published scripts for spectral parameters which probe the position or depth of a given spectral absorption feature. Batch system scalars commonly use a 3-band polynomial fit, while the User Scalars employ a Multiple Feature Extraction Method (Laukamp et al., 2010) for their outputs so are much more restrictive. Details of the scalars name, application, as well as references are included the explanatory notes of the report and in Laukamp et al., 2021.

The TSG[™] file also includes some of the geochemistry data compiled by GSWA (Appendix "Assay_C3DD024_dmirs.csv"), which can be found under the scalars as imported numeric logs and are displayed in plot layout "XRF geochem log". A suite of 37 elements (Al₂O₃, As₂O₃, BaO, Bi₂O₃, CaO, CeO₂, CoO, Cr₂O₃, Cs₂O, CuO, Fe₂O₃, HfO₂, K₂O, La₂O₃, Li₂O, MgO, MnO, Na₂O, Nb₂O₅, NiO, P₂O₅, PbO, Rb₂O, Sb₂O₃, SiO₂, SnO₂, SO₃, SrO, Ta₂O₅, ThO₂, TiO₂, U₃O₈, V₂O₅, WO₃, Y₂O₃, ZnO, ZrO₂) was obtained by GSWA from ~1m intervals of half drill core by means of XRF in 2015.

The lithology data (Appendix "Geology_C3DD024_dmirs2.csv") is captured as class scalars and are displayed in plot layout "lithology log".

Downhole Summary

The most basic hyperspectral output from HyLogger datasets is the downhole mineralogy summary. This is displayed on the summary screen in TSGTM and provides an overview of the downhole hyperspectral mineralogy as derived from the TSA algorithm and can be binned at different sizes. Figure 6 shows the TSA derived downhole mineralogy for C3DD024, the major lithologies and whole rock geochemistry assays Li₂O, K₂O, Na₂O and Sn₂O.

The TSA outputs in Figure 6 are split into SWIR and TIR mineralogy and represent spatial plots of the User SWIR and User jCLST results binned at the 1 m scale. Using only this summary plot, the major lithologies and contacts can be distinguished based on changes in the downhole spectral mineralogy. For instance, the contact between the pegmatite and footwall amphibolite is clearly identifiable by a marked increase in the proportion of amphibole group minerals in the bottom ~10 meters. The crosscutting dolerites are also easily distinguished by amphibole rich sections in the drill core.

It is also clear that while the drill hole intersects over 200 meters of pegmatite, that there are two mineralogical zones which are intersected in the pegmatite based on the thermal-infrared TSA results. The top 120 m of the drill hole contains pyroxene and a strong quartz signature, and the bottom 120 m is characterized by an abundance of K-feldspar, plagioclase and less quartz. This change in pegmatite mineralogy is more subtle in the SWIR results where SWIR TSA results are potentially dominated by smectite in the top 120 m with a notable tourmaline signature below that. However, this does not mean that the bulk mineralogy of the core is dominated by smectites. It is important to consider both the SWIR and TIR together, and which minerals are active in each spectral range when drawing conclusions from hyperspectral mineralogy. Part II of this case study will investigate the smectite and white mica SWIR signature further, in order to determine which di-octahedral sheet silicates are actually present in the respective drill core intervals.

The change in pegmatite mineralogy can also be observed in the core photography (Figure 4) and is also reflected in the downhole geochemistry (Figure 6). The top 120 m comprises the highest Li₂O assay values which reflects the presence of spodumene (Li-pyroxene). The abundance of K-feldspar and plagioclase group minerals (albite) below that spodumene-quartz domain corresponds to higher Na₂O and K₂O values with respect to the top 120 m (Figure 6). This lower Feldspar-plagioclase-tourmaline domain is also characterized by higher SnO₂ values.

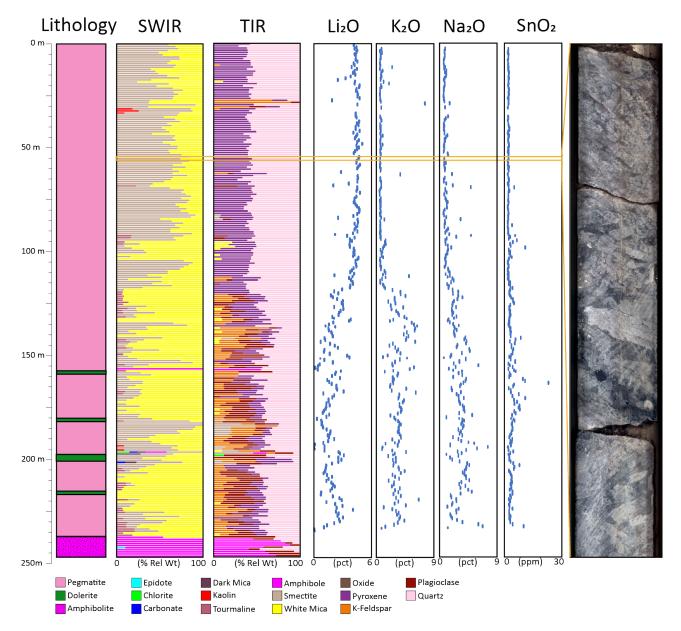


Figure 6 Summary of the major lithologies, TSA-derived hyperspectral mineralogy by mineral group (after Hancock and Wawryk, 2020), and selected goechemical assays for C3DD024, as well as a core photo of a spodumene rich section

Saved plot layouts

A range of plot layouts are saved in the .tdg-file to help navigate the hyperspectral and geochemical data. The plot layouts in TSG can be accessed by selecting "View" > "Plot layouts". For scripts applied in the plot layouts see Laukamp et al. (2021). For details about applying PLS-modelling of geochemical parameters to hyperspectral data see Laukamp et al. (2020).

1.1 Plot layout "log mineralmatch"

This plot layout features from left to right the depth (1st column), user-based TSA SWIR results of the first three mineral groups identified in the SWIR (columns 2, 3 and 4) and the related error (column 5), system-based jCLS TIR results of the first three mineral groups identified in the TIR (columns 6, 7, 8) and the related error (column 9), and the high-resolution linescan RGB imagery (column 10).



Figure 7 Plot layout "log mineralmatch"

1.2 Plot layout "log XRF geochem"

This plot layout features from left to right the whole rock XRF geochemistry results provided by GSWA (Al₂O₃, As₂O₃, BaO, Bi₂O₃, CaO, CeO₂, CoO, Cr₂O₃, Cs₂O, CuO, Fe₂O₃, HfO₂, K₂O, La₂O₃, Li₂O, MgO, MnO, Na₂O, Nb₂O₅, NiO, P₂O₅, PbO, Rb₂O, Sb₂O₃, SiO₂, SnO₂, SO₃, SrO, Ta₂O₅, ThO₂, TiO₂, U₃O₈, V₂O₅, WO₃, Y₂O₃, ZnO, ZrO₂).

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		NULL	NULL	0.061	50	NULL	0.024	NULL	0.075	0.346	4.162	0.043	0.04	0.486	NULL	NULL	0.087	NULL	0.091	39		59	0.016	40	NULL	NULL	NULL	0.045	NULL	NULL	NULL
		NULL	NULL	0.061	50	NULL	0.024	NULL	0.075	0.346	4.182	0.043	0.04	0.486	NULL	NULL	0.087	NULL	0.091	39		59	0.016	48	NULL	NULL	NULL	0.045	NULL	NULL	NULL
		NULL	NULL	0.061 0.061	50 50	NULL	0.024	NULL	0.075	0.346	4.182 4.182	0.043	0.04	0.486	NULL	NULL	0.087	NULL	0.091	39 39		59 59	0.016	48	NULL	NULL	NULL	0.045	NULL	NULL	NULL
		NULL	NULL	0.061	50	NULL	0.024	NULL	0.075	0.346	4.182	0.043	0.04	0.486	NULL	NULL	0.087	NULL	0.091	39		59	0.016	48	NULL	NULL	NULL	0.045	NULL	NULL	NULL
NU	JLL	NULL	NULL	0.061	50	NULL	0.024	NULL	0.075	0.346	4.182	0.043	0.04	0.486	NULL	NULL	0.087	NULL	0.091	39	76.462	59	0.016	48	NULL	NULL	NULL	0.045	NULL	NULL	NULL
		NULL	NULL	0.061	50	NULL	0.024	NULL	0.075	0.346	4.182	0.043	0.04	0.486	NULL	NULL	0.087	NULL	0.091	39		59	0.016	48	NULL	NULL	NULL	0.045	NULL	NULL	NULL
	JLL	NULL	NULL	0.061	50 50	NULL	0.024	NULL	0.075	0.346	4.182	0.043	0.04	0.486	NULL	NULL	0.087	NULL	0.091	39 39		59 59	0.016	48	NULL	NULL	NULL	0.045	NULL	NULL	NULL
NU	JLL	NULL	NULL	0.061	50	NULL	0.024	NULL	0.075	0.346	4.182	0.043	0.04	0.486	NULL	NULL	0.087	NULL	0.091	39	76.462	59	0.016	48	NULL	NULL	NULL	0.045	NULL	NULL	NULL
NL	JLL	NULL	NULL	0.061	50	NULL	0.024	NULL	0.075	0.346	4.182	0.043	0.04	0.486	NULL	NULL	0.087	NULL	0.091	39	76.462	59	0.016	48	NULL	NULL	NULL	0.045	NULL	NULL	NULL
N	AL	NULL	NULL	0.061	50	NULL	0.024	NULL	0.075	0.346	4.162	0.043	0.04	0.486	NULL	NULL	0.087	NULL	0.091	39	16.462	23	0.016	48	NULL	NULL	NULL	0.045	NULL	NULL	NULL

Figure 8 Plot layout "log XRF geochem"

1.3 Plot layout "lithology"

This plot layout features from left to right the depth (1st column), lithological code (2nd column) and the high-resolution linescan RGB imagery (3rd column).

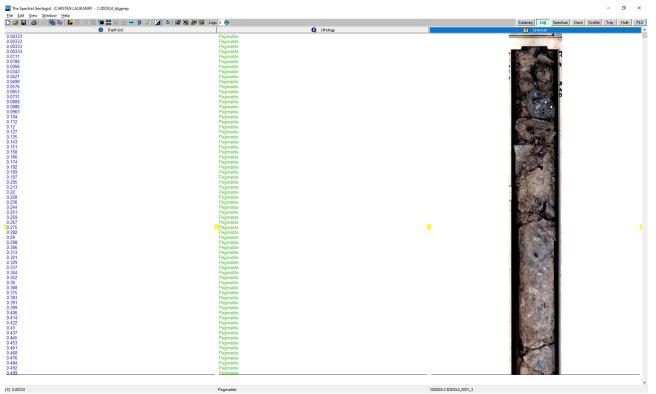


Figure 9 Plot layout "log lithology"

1.4 Plot layout "scatter spodumene"

This plot layout shows VNIR-SWIR reflectance spectra (top left) and TIR reflectance spectra (middle left) as well as the following scatter plots:

- top right: depth (x-axis), XRF-derived Li₂O (y-axis), coloured by the relative intensity of a trough at around 11500 nm diagnostic for spodumene (11500D)
- middle right: depth (x-axis), relative intensity of a trough at around 11500 nm diagnostic for spodumene (11500D) (y-axis), coloured by the XRF-derived Li₂O
- bottom right: depth (x-axis), Aux-match of spodumene reference sample M2082 (y-axis), coloured by the relative intensity of a trough at around 11500 nm diagnostic for spodumene (11500D)
- bottom left: XRF-derived Li₂O (x-axis), relative intensity of a trough at around 11500 nm diagnostic for spodumene (11500D) (y-axis), coloured by the system-based jCLS TIR results of the first mineral group identified in the TIR

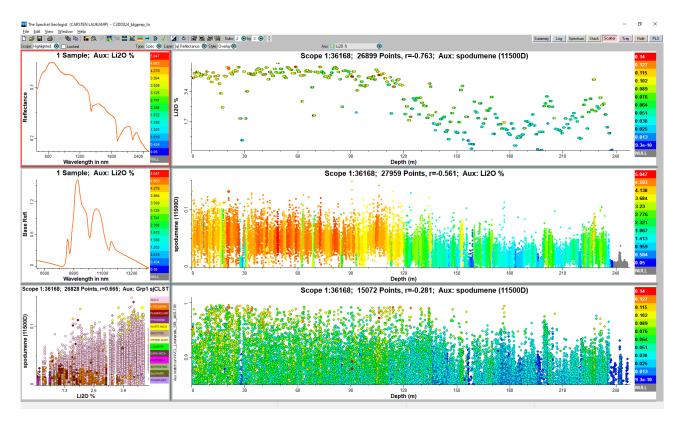


Figure 10 Plot layout "scatter spodumene"

1.5 Plot layout "log PLS"

This plot layout features from left to right the depth (1^{st} column), three pairs of XRF-derived and PLS-modelled geochemical parameters (K_2O , Li_2O , SiO_2 , K_2O+Na_2O/SiO_2) in columns 2 to 9, selection of calibration (C) and test (T) samples used for the PLS-based prediction of the respective geochemical parameters (column 10) and the relative intensity of a trough at around 11500 nm diagnostic for spodumene (11500D) (column 11).

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🚯 Depth					302 %	PLSPied SiD2 %	K20+Na20/Si02	PLSPred_K20+Na			mene (11500D)
3	NULL	NULL	NULL	NULL	76.462	NULL	NULL	NULL	2073/02	NULL	iene (mode)
3	NULL	NULL	NULL	NULL	76.462	NULL	NULL	NULL	Ť	NULL	
ă	NULL	NULL	NULL	NULL	76.462	NULL	NULL	NULL	ċ	NULL	
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3	NULL	NULL	NULL	NULL	76.462	NULL	NULL	NULL	C	NULL	
3	NULL	NULL	NULL	NULL	76.462	NULL	NULL	NULL	I.	NULL	
	NULL 0.346	NULL 1.603	NULL 4.182	NULL 0.681	76.462 76.462	NULL 73.318	NULL 0.0109	NULL 0.0558	ç	NULL 0.0385	
	0.346	2,193	4,182	0.71	76.462	72.266	0.0109	0.0746		0.0352	
	0.346	2.632	4,182	0.919	76.462	71.975	0.0109	0.0673	Ť	0.0492	
	0.346	3.179	4,182	0.778	76.462	71.559	0.0109	0.0753	Ċ	0.038	
	0.346	2.856	4.182	0.628	76.462	71.683	0.0109	0.075	т	0.0448	
	0.346	3.137	4.182	0.64	76.462	71.959	0.0109	0.0665	ç	0.0353	
	0.346	NULL	4.182	NULL	76.462	NULL	0.0109	NULL	T	0.00463	
	0.346 0.346	NULL	4.182 4.182	NULL	76.462 76.462	NULL	0.0109 0.0109	NULL	ç	0.00464	
	0.346	NULL	4.182	NULL	76.462	NULL	0.0109	NULL	i i	0.00669	
	0.346	NULL	4.182	NULL	76.462	NULL	0.0109	NULL	Ť	0.00302	
	0.346	NULL	4.182	0.688	76.462	71.247	0.0109	0.0886	ċ	0.0214	
	0.346	1.538	4.182	1.117	76.462	73.55	0.0109	0.0368	т	0.0629	
	0.346	2.334	4.182	1.321	76.462	NULL	0.0109	NULL	С	0.0677	
	0.346	NULL	4.182	NULL	76.462	NULL	0.0109	NULL	т	0.091	
	0.346	NULL	4.182	NULL	76.462	NULL	0.0109	NULL	ç	0.0619	
	0.346 0.346	NULL	4.182 4.182	NULL	76.462 76.462	NULL	0.0109 0.0109	NULL	L L	0.132 0.163	
	0.346	NULL	4.102	1.953	76.462	NULL	0.0109	NULL	¥	0.163	
	0.346	3.732	4.182	0.6	76.462	71.282	0.0109	0.0895		0.0213	
	0.346	4.098	4.182	0.6	76.462	70.798	0.0109	0.0939	Ť	0.0075	
	0.346	3.583	4.182	0.562	76.462	71.44	0.0109	0.0895	ċ	0.0075	
	0.346	2.684	4.182	0.715	76.462	72.707	0.0109	0.0742	Ť	0.0312	
	0.346	2.713	4.182	0.584	76.462	NULL	0.0109	0.0691	С	0.0288	
	0.346	2.013	4.182	0.639	76.462	NULL	0.0109	NULL	Ţ	0.0185	
	0.346	NULL	4.182	0.707	76.462	NULL	0.0109	NULL	ç	0.0331	
	0.346 0.346	NULL 2.304	4.182 4.182	0.693 0.806	76.462 76.462	NULL 73.194	0.0109 0.0109	NULL 0.0622	L L	0.0195 0.0478	
	0.346	2.454	4.162	1.182	76.462	73.287	0.0109	0.0506	÷ ÷	0.0478	
	0.346	2.429	4,182	0.834	76.462	72.725	0.0109	0.0645	i i	0.0319	
	0.346	2.025	4,182	0.752	76.462	72.613	0.0109	0.0574	Ť	0.0354	
	0.346	3.034	4.182	1.043	76.462	72.436	0.0109	0.0675	ċ	0.067	
	0.346	2.418	4.182	0.84	76.462	72.068	0.0109	0.0588	т	0.0354	
	0.346	2.855	4.182	0.949	76.462	71.461	0.0109	0.0604	С	0.0294	
	0.346	2.533	4.182	0.889	76.462	72.296	0.0109	0.0624	T. T.	0.0313	
	0.346 0.346	1.585	4.182 4.182	NULL	76.462 76.462	73.435 73.708	0.0109 0.0109	0.0359 0.0303	ç	0.0389 0.0451	
	0.346	1.242 2.477	4.182	1.275	76.462	73.708	0.0109	0.0303	L.	0.0538	
	0.346	2.311	4.182	0.989	76.462	72.211	0.0109	0.0528	Ť	0.0415	
	0.346	1.847	4.182	1.112	76.462	73.054	0.0109	0.0429	ċ	0.0489	
	0.346	2.528	4.182	1.673	76.462	73.609	0.0109	0.0507	Ť	0.0641	
	0.346	2.084	4.182	1.046	76.462	73.373	0.0109	0.0576	Ċ	0.049	
	0.346	2.069	4.182	1.283	76.462	73.514	0.0109	0.0552	Ţ	0.0534	
	0.346	2.06	4.182	1.232	76.462	73.489	0.0109	0.051	ç	0.0466	
	0.346	2.321 2.238	4.182	1.111	76.462	72.635	0.0109	0.0515	T	0.0466	
	0.346	2.238	4.182 4.182	1.633 NULL	76.462 76.462	72.987 73.413	0.0109	0.0493 0.0502	ç	0.065 0.0725	
	0.346	1.935	4.102	1.599	76.462	73.793	0.0109	0.0502	i.	0.0725	
	0.346	1.253	4,182	1.272	76.462	74.264	0.0109	0.0498	Ť	0.0559	
	0.346	1.787	4.182	1.133	76.462	73.619	0.0109	0.0513	ċ	0.0412	
	0.346	2.512	4.182	0.794	76.462	72.596	0.0109	0.0631	Ť	0.034	
	0.346	2.691	4.182	0.146	76.462	72.492	0.0109	0.0673	С	0.005	
	0.346	2.287	4.182	0.59	76.462	72.429	0.0109	0.0561	Ţ	0.0218	
	0.346	1.56	4.182	0.902	76.462	73.549	0.0109	0.045	ç	0.0414	
	0.346	1.766	4.182 4.182	0.865	76.462	73.662	0.0109	0.0485	1	0.0406 0.0527	
	0.346 0.346	2.047 2.777	4.182	1.108	76.462 76.462	72.721 71.861	0.0109 0.0109	0.0616 0.0647	÷ ÷	0.0527	
	0.346	2.225	4.182	0.881	76.462	72.217	0.0109	0.0653		0.0523	
	0.346	2.294	4.182	0.964	76.462	72.89	0.0109	0.0576	Ť	0.0472	
	0.346	2.411	4.182	0.887	76.462	72.196	0.0109	0.0741	ċ	0.0402	
	0.346	2.803	4.182	0.313	76.462	72.554	0.0109	0.0649	T	0.0111	
133	NULL	NULL	NULL	NULL	76.462	NULL	NULL	NULL	c	NULL	

Figure 11 Plot layout "log PLS"

1.6 Plot layout "scatter PLS K"

This plot layout comprises scatter plots:

- Top left: XRF-derived SiO₂ (x-axis), XRF-derived K₂O (y-axis), coloured by XRF-derived Li₂O
- Top right: depth (x-axis), XRF-derived K₂O (y-axis), selection of calibration (C) and test (T) samples used for the PLS-based prediction of the respective geochemical parameters
- Middle left: XRF-derived K₂O (x-axis), PLS-based modelling of K₂O using only the 50 to 250 m depth interval (y-axis), coloured by the system-based jCLS TIR results of the first mineral group identified in the TIR
- Middle right: depth (x-axis), PLS-based modelling of K₂O using only the 50 to 250 m depth interval (y-axis), coloured by XRF-derived K₂O
- Bottom left: TIR reflectance spectra
- Bottom right: depth (x-axis), XRF-derived K₂O (y-axis), coloured by PLS-based modelling of K₂O using only the 50 to 250 m depth interval

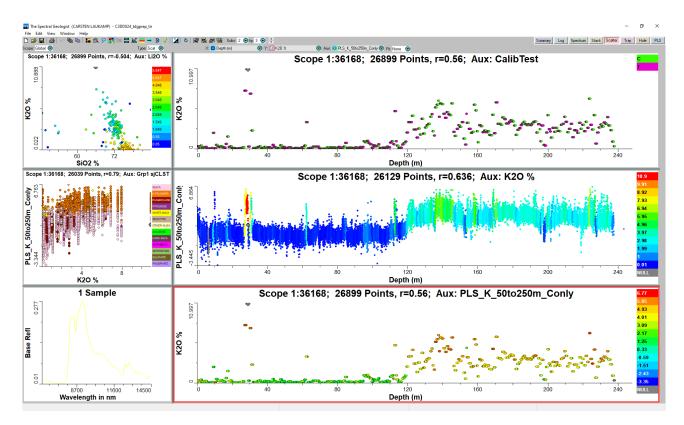


Figure 12 Plot layout "scatter PLS K"

1.7 Plot layout "scatter PLS KNaSi"

This plot layout comprises scatter plots:

- Top left: XRF-derived LiO₂ (x-axis), K₂O+Na₂O/SiO₂ (y-axis), coloured by XRF-derived Li₂O
- Top right: depth (x-axis), K₂O+Na₂O/SiO₂ (y-axis), selection of calibration (C) and test (T) samples used for the PLS-based prediction of the respective geochemical parameters
- Middle left: K₂O+Na₂O/SiO₂ (x-axis), PLS-based modelling of K₂O+Na₂O/SiO₂ (y-axis), coloured by the system-based jCLS TIR results of the first mineral group identified in the TIR
- Middle right: depth (x-axis), PLS-based modelling of K₂O+Na₂O/SiO₂ (y-axis), coloured by K₂O+Na₂O/SiO₂
- Bottom left: TIR reflectance spectra
- Bottom right: depth (x-axis), K₂O+Na₂O/SiO₂ (y-axis), coloured by PLS-based modelling of K₂O+Na₂O/SiO₂

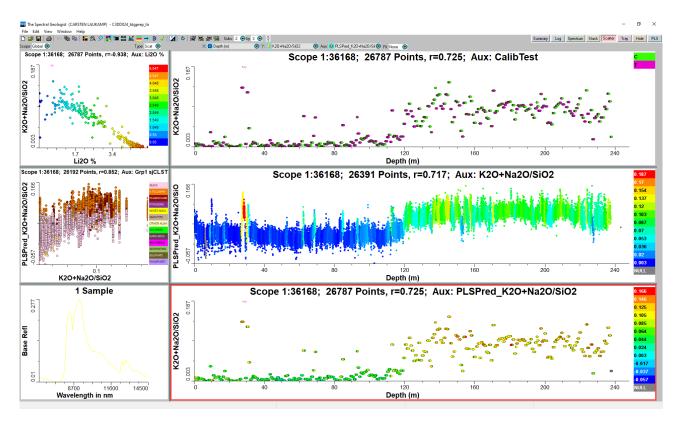


Figure 13 Plot layout "scatter PLS KNaSi"

1.8 Plot layout "tray8 KNaSi PLS"

This plot layout shows the high resolution RGB imagery of tray 8. The top colour ribbon of each section shows K_2O+Na_2O/SiO_2 calculated from XRF results. The bottom colour ribbon shows the PLS-based K_2O+Na_2O/SiO_2 .

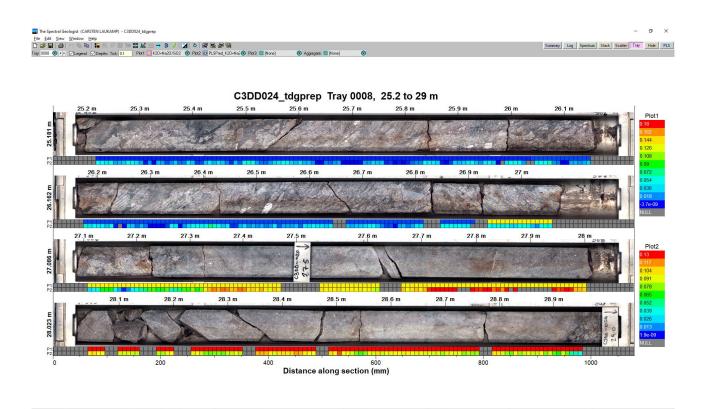


Figure 14 Plot layout "tray8 KNaSi PLS"

Acknowledgments

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Furthermore, we'd like to acknowledge all those hard-working people that collect and process the NVCL data on a day to day basis, helping to 'uncover the mineralogy of the top two kilometres of the Australian continent', and particular, the HyLogging team at the Geological Survey of Western Australia for creating this dataset.







Geological Survey of Western Australia

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GSWA WAMEX Report A0104943

https://www.talisonlithium.com/greenbushes-project

Additional Resources

AuScope Discovery Portal http://portal.auscope.org/

Instructional Videos

"How to download TSG Files from the Auscope Discover Portal" https://www.youtube.com/watch?v=Oh_YAxEpLeo&t=115s

"Introduction to working with HyLogger Data in The Spectral Geologist (TSG)" https://www.youtube.com/watch?v=u-SjA2_J3RQ&t=71s

Recorded Webinars

"The National Virtual Core Library – Building a Continental-Scale drill core database" https://www.youtube.com/watch?v=AlcIqAGc9U4&t=4s

"Mineral composition trends in hydrothermal mineral systems inferred from reflectance spectra" https://www.youtube.com/watch?v=IEvxn6o0imk&t=110s

Other TSG Materials

https://research.csiro.au/thespectralgeologist/support/downloads/

CSIRO Spectral Reference Library

http://mineralspectrallibraries.csiro.au/

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