

# RUNNING THE BLM SOFTWARE

April 6, 2017

ACWA Workshop on Oregon's Copper Standard  
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Water Resources Analyst

Clean Water  Services

## Goals of this Portion of the Workshop

- Show how we have been using the model to generate WQ Criteria for specific sampling events
- Provide tips for getting large datasets ready to input into the model software, including estimating missing parameters needed to run the model
- Step-by-step instructions for running the model
- Show an example dataset used to run the model
- Display results, showing the WQ Criteria generated for the example dataset



# User's Guide by Windward Environmental

- **BIOTIC LIGAND MODEL WINDOWS® INTERFACE,  
RESEARCH VERSION 3.1.2.37: USER'S GUIDE AND  
REFERENCE MANUAL**

- October 2015



# BLM Interface

BLM Freshwater version 3.1.2.37

File Edit Inputs Help

Icons: [New] [Open] [Save] [Print] [Cut] [Copy] [Paste] [Undo] [Redo] [Checkmark] [Help]

Description:

Current Selections

Prediction Mode: Toxicity      Metal: Cu      Organism/Type: US EPA WQC calculation

Site Chemistry    Simplified Site Chemistry

	Site Name	Sample Name	Temp.	pH	Cu	DOC	HA	Ca	Mg	Na	K	SO4	Cl	Alkalinity	S
			C		ug/L	mg C/L	%	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L CaCO3	mg/L
1															
2															
3															
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# Example Dataset – Parameters Missing

Sample Point Name	Calendar Date	00010 - Temperature	00400 - pH	01040 - Cu-Sol	88998 - NPOC	HA	00916 - Ca	00927 - Mg	00929 - Na	00937 - K	00946 - SO4	00941 - Cl	00410 - Alkalinity	S
		°C	S.U.	µg/L	mg/L	%	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L
TR@FERN HILL	1/19/2016	7.9	7.3										3.29	22.6
TR@FERN HILL	2/9/2016	7.7	7.15	0.619	1.17		6.81	2.35	4.07	0.538	2.95	3.39	28.5	
TR@FERN HILL	2/23/2016	6.6	7.41									3.35	29.2	
TR@FERN HILL	3/1/2016	8	7.01	0.792	1.21		7.07	2.54	4.28	0.572	2.93	3.32	29	
TR@FERN HILL	3/8/2016	7.9	7.2									2.88	27	
TR@FERN HILL	3/22/2016	9	7.15									3.19	29.2	
TR@FERN HILL	4/5/2016	10.5	7.19									3.55	35	
TR@FERN HILL	4/13/2016	10.9	7.12		1.43							3.85		
TR@FERN HILL	4/19/2016	15.2	7.21									4.18	36.8	
TR@FERN HILL	4/26/2016	11.3	7.25		1.43							4.42		
TR@FERN HILL	5/2/2016	14.3	7.23									4.39	41.2	
TR@FERN HILL	5/16/2016	12	7.42	0.715	1.21		8.3	2.72				3.96	36.1	
TR@FERN HILL	6/13/2016	13.2	7.49	0.694	1.2		7.41	2.38	4.33	0.447	2.12	3.65	33	
TR@FERN HILL	6/27/2016	14.2	7.42									3.52	29.7	
TR@FERN HILL	7/11/2016	12.3	7.57	0.661	1.19		6.16	2.15	3.66	0.37	2.12	3.57	29.2	
TR@FERN HILL	7/25/2016	11.8	7.27									3.09	27.8	
TR@FERN HILL	8/15/2016	12.4	7.46		1.09						2.09	2.92	26.9	



# DEQ's Specific Conductance Regression Equations

Uses Specific Conductance (X) to estimate the Missing Parameter (Y)

Parameter	Regression Equation	Adjusted R <sup>2</sup>	p-value
DOC	$\ln(y) = 0.69 \cdot \ln(x) - 2.43$	0.31	< 0.001
Hardness	$\ln(y) = 1.02 \cdot \ln(x) - 1.16$	0.92	< 0.001
Alkalinity	$\ln(y) = 0.88 \cdot \ln(x) - 0.41$	0.77	< 0.001
Calcium	$\ln(y) = 0.96 \cdot \ln(x) - 2.29$	0.89	< 0.001
Sodium	$\ln(y) = 0.86 \cdot \ln(x) - 2.22$	0.82	< 0.001
Magnesium	$\ln(y) = 0.91 \cdot \ln(x) - 3.09$	0.85	< 0.001
Potassium	$\ln(y) = 0.84 \cdot \ln(x) - 3.74$	0.70	< 0.001
Chloride	$\ln(y) = 0.15 \cdot \ln(x) - 3.82$	0.77	< 0.001
Sulfate	$\ln(y) = 1.45 \cdot \ln(x) - 5.59$	0.76	< 0.001



# DEQ's Specific Conductance Regression Equations

Parameter	Equation
Alkalinity	$\text{Alk.} = \exp^{(0.88 \cdot [\ln(\text{SpC})] - 0.41)}$
Calcium	$\text{Ca} = \exp^{(0.96 \cdot [\ln(\text{SpC})] - 2.29)}$
Chloride	$\text{Cl} = \exp^{(1.15 \cdot [\ln(\text{SpC})] - 3.82)}$
Magnesium	$\text{Mg} = \exp^{(0.91 \cdot [\ln(\text{SpC})] - 3.09)}$
Potassium	$\text{K} = \exp^{(0.84 \cdot [\ln(\text{SpC})] - 3.74)}$
Sodium	$\text{Na} = \exp^{(0.86 \cdot [\ln(\text{SpC})] - 2.22)}$
Sulfate	$\text{SO}_4 = \exp^{(1.45 \cdot [\ln(\text{SpC})] - 5.59)}$



# Example Dataset – Missing Parameters Estimated (Except Humic Acid and Sulfide)

Sample Point Name	Calendar Date	00010 - Temperature	00400 - pH	01040 - Cu-Sol	88998 - NPOC	HA	00916 - Ca	00927 - Mg	00929 - Na	00937 - K	00946 - SO4	00941 - Cl	00410 - Alkalinity	S
		°C	S.U.	µg/L	mg/L	%	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L
TR@FERN HILL	1/19/2016	7.9	7.3			1.65	6.42	2.23	3.73	0.518	2.45	3.29		22.6
TR@FERN HILL	2/9/2016	7.7	7.15	0.619		1.17	6.81	2.35	4.07	0.538	2.95	3.39		28.5
TR@FERN HILL	2/23/2016	6.6	7.41			1.31	7.10	2.42	4.12	0.676	2.87	3.35		29.2
TR@FERN HILL	3/1/2016	8	7.01	0.792		1.21	7.07	2.54	4.28	0.572	2.93	3.32		29
TR@FERN HILL	3/8/2016	7.9	7.2			1.33	6.52	2.26	3.79	0.541	2.51	2.88		27
TR@FERN HILL	3/22/2016	9	7.15			1.44	7.30	2.47	4.23	0.723	2.99	3.19		29.2
TR@FERN HILL	4/5/2016	10.5	7.19			1.77	8.00	2.66	4.63	0.884	3.42	3.55		35
TR@FERN HILL	4/13/2016	10.9	7.12			1.43	8.32	2.75	4.81	0.958	3.61	3.85		34.1
TR@FERN HILL	4/19/2016	15.2	7.21			2.74	8.53	2.81	4.93	1.007	3.74	4.18		36.8
TR@FERN HILL	4/26/2016	11.3	7.25			1.43	8.83	2.89	5.11	1.077	3.92	4.42		36.1
TR@FERN HILL	5/2/2016	14.3	7.23			2.62	9.35	3.03	5.40	1.196	4.24	4.39		41.2
TR@FERN HILL	5/16/2016	12	7.42	0.715		1.21	8.3	2.72	4.74	0.927	3.53	3.96		36.1
TR@FERN HILL	6/13/2016	13.2	7.49	0.694		1.2	7.41	2.38	4.33	0.447	2.12	3.65		33
TR@FERN HILL	6/27/2016	14.2	7.42			1.19	6.98	2.38	4.05	0.647	2.79	3.52		29.7
TR@FERN HILL	7/11/2016	12.3	7.57	0.661		1.19	6.16	2.15	3.66	0.37	2.12	3.57		29.2
TR@FERN HILL	7/25/2016	11.8	7.27			1.11	6.52	2.26	3.79	0.541	2.51	3.09		27.8
TR@FERN HILL	8/15/2016	12.4	7.46			1.09	6.43	2.23	3.74	0.520	2.09	2.92		26.9





# BLM Interface

BLM Freshwater version 3.1.2.37

File Edit Inputs Help

Icons: [New] [Open] [Save] [Print] [Cut] [Copy] [Paste] [Undo] [Redo] [Checkmark] [Help]

Description:

Current Selections

Prediction Mode: Toxicity      Metal: Cu      Organism/Type: US EPA WQC calculation

Site Chemistry    Simplified Site Chemistry

	Site Name	Sample Name	Temp.	pH	Cu	DOC	HA	Ca	Mg	Na	K	SO4	Cl	Alkalinity	S
			C		ug/L	mg C/L	%	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L CaCO3	mg/L
1															
2															
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# BLM Interface

BLM Freshwater version 3.1.2.37

File Edit Inputs Help

Description:

Current Selections  
Prediction Mode: Toxicity      Metal: Cu      Organism/Type: US EPA WQC calculation

Site Chemistry    Simplified Site Chemistry

	Site Name	Sample Name	Temp.	pH	Cu	DOC	HA	Ca	Mg	Na	K	SO4	Cl	Alkalinity	S
			C		ug/L	mg C/L	%	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L CaCO3	mg/L
1															
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Note the order of the parameters in the columns.  
If you plan to run the model for a large number of samples simultaneously, using this order in your spreadsheet will make it easy to paste into the BLM interface



# Column Order of the Parameters

## Order Parameter

- 1 Site Name
- 2 Sample Name
- 3 Temperature
- 4 pH
- 5 Copper (not necessary to run the model)
- 6 Dissolved Organic Carbon
- 7 Humic Acid
- 8 Calcium
- 9 Magnesium
- 10 Sodium
- 11 Potassium
- 12 Sulfate
- 13 Chloride
- 14 Alkalinity
- 15 Sulfide



# BLM Interface

BLM Freshwater version 3.1.2.37

File Edit Inputs Help

Description: |

Current Selections  
Prediction Mode: Toxicity      Metal: Cu      Organism/Type: US EPA WQC calculation

Site Chemistry    Simplified Site Chemistry

	Site Name	Sample Name	Temp.	pH	Cu	DOC	HA	Ca	Mg	Na	K	SO4	Cl	Alkalinity	S
			C		ug/L	mg C/L	%	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L CaCO3	mg/L
1															
2															
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Note the units used for each parameter. Units displayed above are defaults for the software, but they can be changed by the user



# BLM Interface

BLM Freshwater version 3.1.2.37

File Edit Inputs Help

Description: |

Current Selections

Prediction Mode: Toxicity Metal: Cu Organism/Type: US EPA WQC calculation

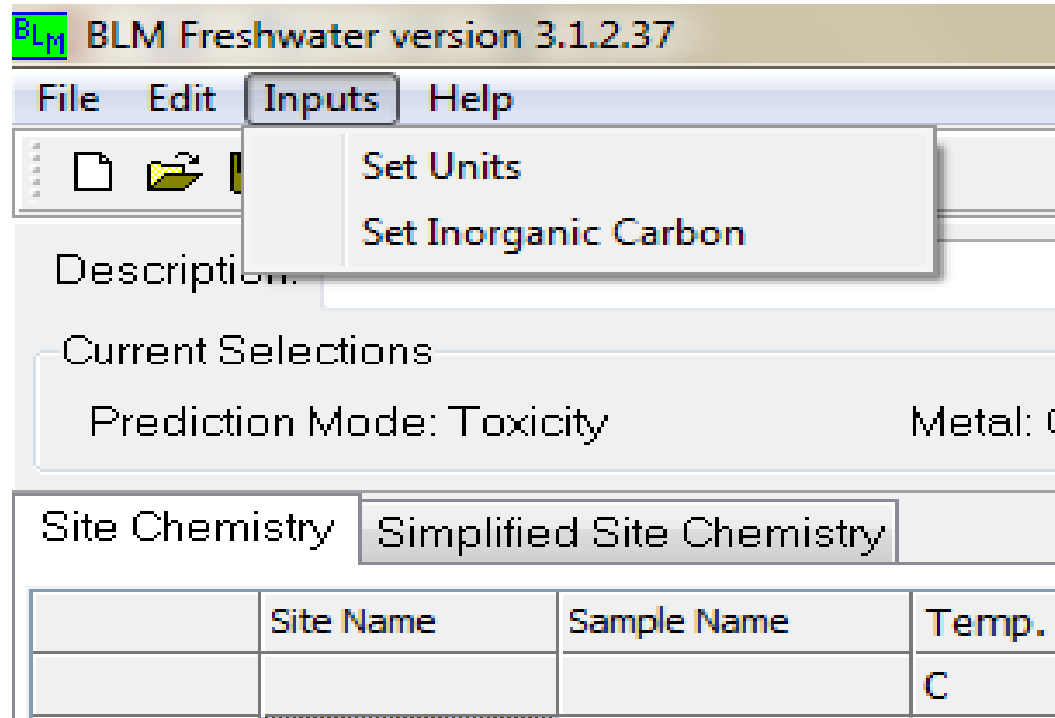
Site Chemistry Simplified Site Chemistry

	Site Name	Sample Name	Temp.	pH	Cu	DOC	HA	Ca	Mg	Na	K	SO4	Cl	Alkalinity	S
1			C		ug/L	mg C/L	%	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L CaCO3	mg/L
2															
3															
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To set units, click the **Inputs** menu item



# Setting the Units

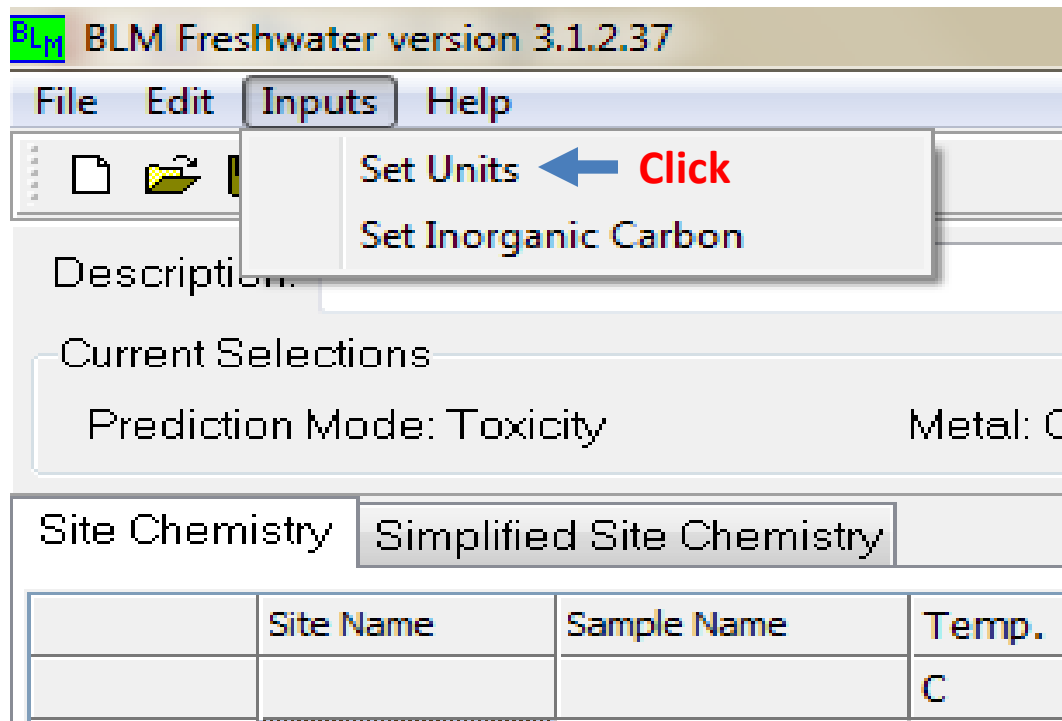


The screenshot shows the BLM Freshwater software interface. The title bar reads "BLM Freshwater version 3.1.2.37". The menu bar includes "File", "Edit", "Inputs", and "Help". The "Inputs" menu is open, showing two options: "Set Units" and "Set Inorganic Carbon". Below the menu, the "Description" field is visible. The "Current Selections" section shows "Prediction Mode: Toxicity" and "Metal: C". The "Site Chemistry" section has "Simplified Site Chemistry" selected. At the bottom, a table displays site information.

	Site Name	Sample Name	Temp.
			C



# Setting the Units

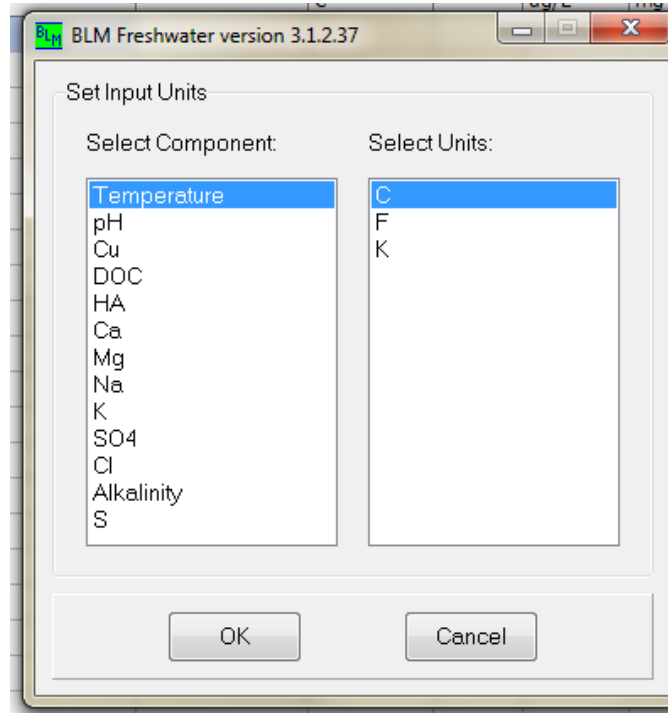


The screenshot shows the BLM Freshwater version 3.1.2.37 application window. The 'Inputs' menu is open, and the 'Set Units' option is highlighted with a blue arrow and the word 'Click' in red. The 'Set Inorganic Carbon' option is also visible. Below the menu, the 'Current Selections' section shows 'Prediction Mode: Toxicity' and 'Metal: C'. The 'Site Chemistry' section shows 'Simplified Site Chemistry' selected. At the bottom, a table displays site and sample information.

	Site Name	Sample Name	Temp.
			C



# Dialogue Box for Setting Parameter Units





# BLM Interface

BLM Freshwater version 3.1.2.37

File Edit Inputs Help

Description: |

Current Selections

Prediction Mode: Toxicity Metal: Cu Organism/Type: US EPA WQC calculation

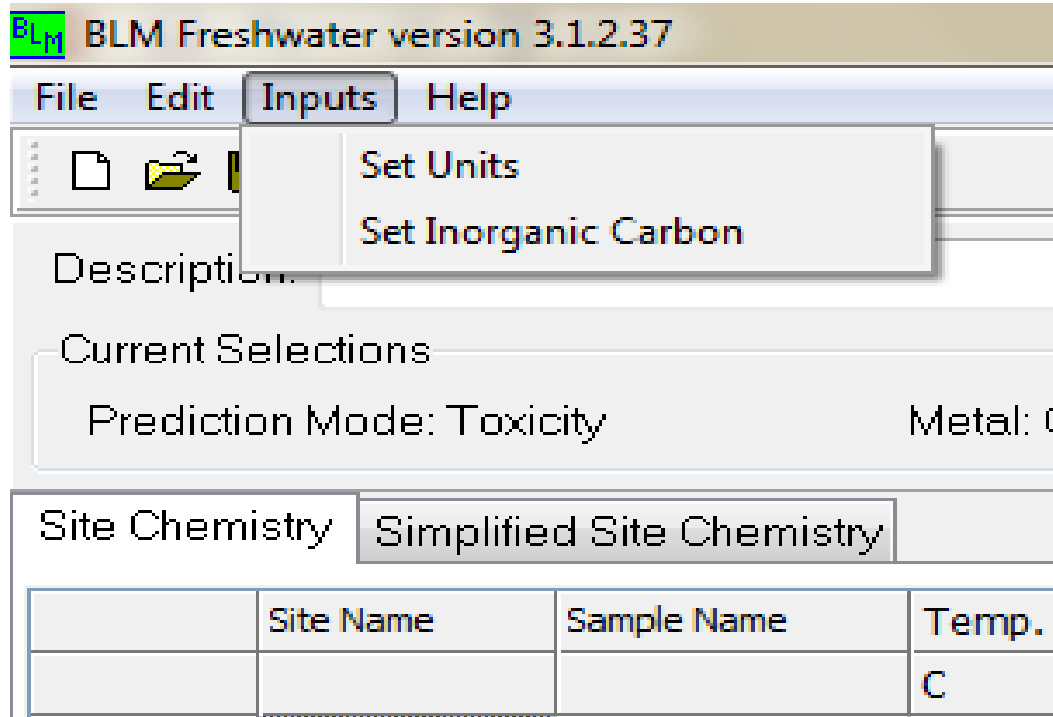
Site Chemistry Simplified Site Chemistry

	Site Name	Sample Name	Temp.	pH	Cu	DOC	HA	Ca	Mg	Na	K	SO4	Cl	Alkalinity	S
			C		ug/L	mg C/L	%	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L CaCO3	mg/L
1															
2															
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To set inorganic carbon input, click the **Inputs** menu item



# Setting Inorganic Carbon Input

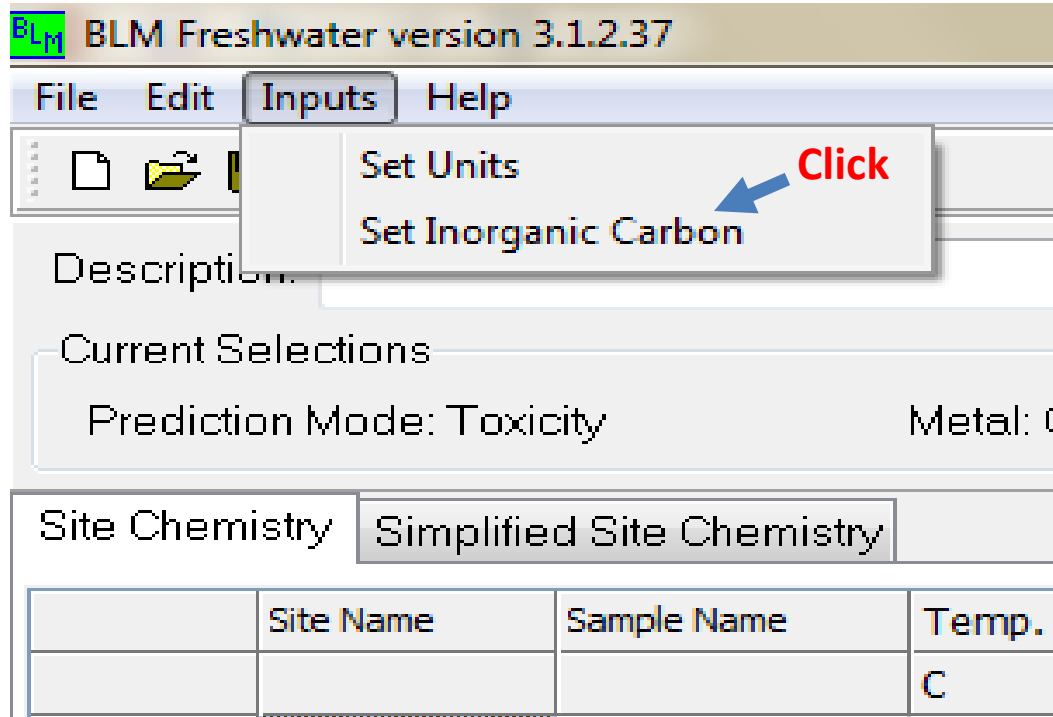


The screenshot shows the BLM Freshwater version 3.1.2.37 application window. The 'Inputs' menu is open, displaying options for 'Set Units' and 'Set Inorganic Carbon'. The 'Set Inorganic Carbon' option is highlighted. Below the menu, the 'Current Selections' section shows 'Prediction Mode: Toxicity' and 'Metal: C'. The 'Site Chemistry' section shows 'Simplified Site Chemistry' selected. At the bottom, a table displays site information.

	Site Name	Sample Name	Temp.
			C



# Setting Inorganic Carbon Input

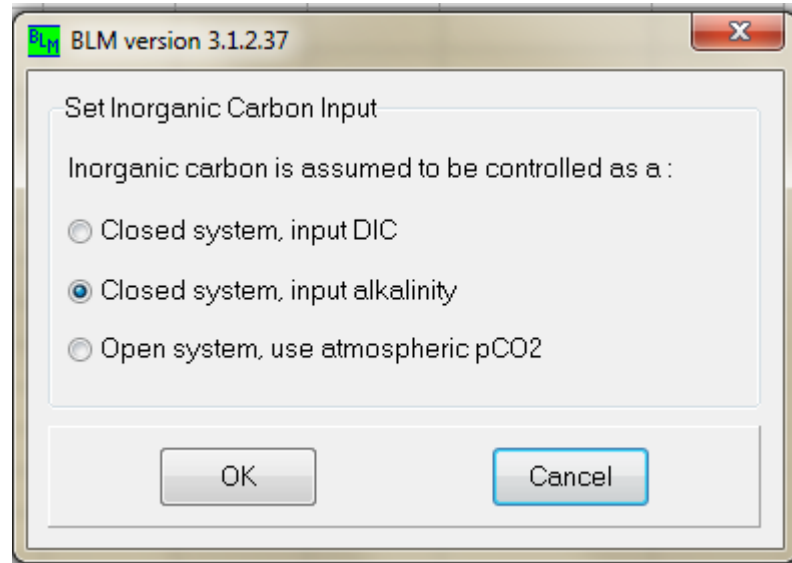


The screenshot shows the BLM Freshwater version 3.1.2.37 application window. The 'Inputs' menu is open, and the 'Set Inorganic Carbon' option is highlighted with a blue arrow and the word 'Click' in red text. The interface includes a menu bar with 'File', 'Edit', 'Inputs', and 'Help'. Below the menu bar, there are icons for file operations. The main area contains a 'Description' field, a 'Current Selections' section with 'Prediction Mode: Toxicity' and 'Metal: C', and a 'Site Chemistry' section with 'Simplified Site Chemistry' selected. At the bottom, there is a table with columns for 'Site Name', 'Sample Name', and 'Temp.', with the value 'C' entered in the 'Temp.' column.

	Site Name	Sample Name	Temp.
			C



# Dialogue Box for Setting Organic Carbon Input



# BLM Interface

BLM Freshwater version 3.1.2.37

File Edit Inputs Help

Description:

Current Selections  
Prediction Mode: Toxicity      Metal: Cu      Organism/Type: US EPA WQC calculation

Site Chemistry    Simplified Site Chemistry

	Site Name	Sample Name	Temp.	pH	Cu	DOC	HA	Ca	Mg	Na	K	SO4	Cl	Alkalinity	S
			C		ug/L	mg C/L	%	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L CaCO3	mg/L
1															
2															
3															
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17															
18															
19															

Site Name: Information about the sample site, e.g., Tualatin River at Fernhill Road



# BLM Interface

BLM Freshwater version 3.1.2.37

File Edit Inputs Help

Description: |

Current Selections  
Prediction Mode: Toxicity      Metal: Cu      Organism/Type: US EPA WQC calculation

Site Chemistry    Simplified Site Chemistry

	Site Name	Sample Name	Temp.	pH	Cu	DOC	HA	Ca	Mg	Na	K	SO4	Cl	Alkalinity	S
			°C		ug/L	mg C/L	%	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L	mg/L CaCO3	mg/L
1															
2															
3															
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19															

Sample Name:  
Information about the  
sample; e.g., the sample  
collection date



# Default Value of 10% for Humic Acid

BLM Freshwater version 3.1.2.37 - H:\My Documents\BLM\BLM software\Forest Grove upstream.blm

File Edit Inputs Help

Description: Forest Grove upstream

Current Selections  
 Prediction Mode: Toxicity      Metal: Cu      Organism/Type: US EPA WQC calculation

Site Chemistry    Simplified Site Chemistry

	Site Name	Sample Name	Temp. C	pH	Cu ug/L	DOC mg C/L	HA %	Ca mg/L	Mg mg/L	Na mg/L	K mg/L	SO4 mg/L	Cl mg/L	Alkalinity mg/L CaCO3	S mg/L
149	TR@FERN HILL	1/19/2016	7.90	7.30	-6.35E+	1.51	10.00	6.42	2.22	3.32	0.52	3.55	3.29	22.60	1.00E-001
150	TR@FERN HILL	2/9/2016	7.70	7.15	0.62	1.17	10.00	6.81	2.35	4.07	0.54	2.95	3.39	28.50	1.00E-001
151	TR@FERN HILL	2/23/2016	6.60	7.41	-6.35E+	1.22	10.00	7.10	2.41	3.73	0.68	3.99	3.35	29.20	1.00E-001
152	TR@FERN HILL	3/1/2016	8.00	7.01	0.79	1.21	10.00	7.07	2.54	4.28	0.57	2.93	3.32	29.00	1.00E-001
153	TR@FERN HILL	3/8/2016	7.90	7.20	-6.35E+	1.24	10.00	6.52	2.25	3.38	0.54	3.62	2.88	27.00	1.00E-001
154	TR@FERN HILL	3/22/2016	9.00	7.15	-6.35E+	1.33	10.00	7.30	2.47	3.85	0.72	4.11	3.19	29.20	1.00E-001
155	TR@FERN HILL	4/5/2016	10.50	7.19	-6.35E+	1.62	10.00	8.00	2.66	4.26	0.88	4.56	3.55	35.00	1.00E-001
156	TR@FERN HILL	4/13/2016	10.90	7.12	-6.35E+	1.40	10.00	8.32	2.75	4.45	0.96	4.76	3.85	34.20	1.00E-001
157	TR@FERN HILL	4/19/2016	15.20	7.21	-6.35E+	2.45	10.00	8.53	2.81	4.58	1.01	4.90	4.18	36.80	1.00E-001
158	TR@FERN HILL	4/26/2016	11.30	7.25	-6.35E+	1.40	10.00	8.84	2.89	4.76	1.08	5.09	4.42	36.10	1.00E-001
159	TR@FERN HILL	5/2/2016	14.30	7.23	-6.35E+	2.34	10.00	9.35	3.04	5.07	1.20	5.42	4.39	41.20	1.00E-001
160	TR@FERN HILL	5/16/2016	12.00	7.42	0.71	1.21	10.00	8.30	2.72	4.37	0.93	4.68	3.96	36.10	1.00E-001
161	TR@FERN HILL	6/13/2016	13.20	7.49	0.69	1.20	10.00	7.41	2.38	4.33	0.45	2.12	3.65	33.00	1.00E-001
162	TR@FERN HILL	6/27/2016	14.20	7.42	-6.35E+	1.12	10.00	6.98	2.38	3.65	0.65	3.90	3.52	29.70	1.00E-001
163	TR@FERN HILL	7/11/2016	12.30	7.57	0.66	1.19	10.00	6.16	2.15	3.66	0.37	2.12	3.57	29.20	1.00E-001
164	TR@FERN HILL	7/25/2016	11.80	7.27	-6.35E+	1.05	10.00	6.52	2.25	3.38	0.54	3.62	3.09	27.80	1.00E-001
165	TR@FERN HILL	8/15/2016	12.40	7.46	0.79	1.09	10.00	6.45	2.08	3.54	0.43	2.09	2.92	26.90	1.00E-001
166	TR@FERN HILL	8/29/2016	14.40	7.50	-6.35E+	1.12	10.00	6.48	2.24	3.36	0.53	3.59	3.00	27.10	1.00E-001
167	TR@FERN HILL	9/12/2016	15.70	7.50	-6.35E+	1.16	10.00	6.76	2.32	3.52	0.60	1.89	3.33	28.60	1.00E-001
168															



# Arbitrary Value for Sulfide Needed to Run the Model

BLM Freshwater version 3.1.2.37 - H:\My Documents\BLM\BLM software\Forest Grove upstream.blm

File Edit Inputs Help

Description: Forest Grove upstream

Current Selections  
 Prediction Mode: Toxicity      Metal: Cu      Organism/Type: US EPA WQC calculation

Site Chemistry    Simplified Site Chemistry

	Site Name	Sample Name	Temp. C	pH	Cu ug/L	DOC mg C/L	HA %	Ca mg/L	Mg mg/L	Na mg/L	K mg/L	SO4 mg/L	Cl mg/L	Alkalinity mg/L CaCO3	S mg/L
149	TR@FERN HILL	1/19/2016	7.90	7.30	-6.35E+	1.51	10.00	6.42	2.22	3.32	0.52	3.55	3.29	22.60	1.00E-001
150	TR@FERN HILL	2/9/2016	7.70	7.15	0.62	1.17	10.00	6.81	2.35	4.07	0.54	2.95	3.39	28.50	1.00E-001
151	TR@FERN HILL	2/23/2016	6.60	7.41	-6.35E+	1.22	10.00	7.10	2.41	3.73	0.68	3.99	3.35	29.20	1.00E-001
152	TR@FERN HILL	3/1/2016	8.00	7.01	0.79	1.21	10.00	7.07	2.54	4.28	0.57	2.93	3.32	29.00	1.00E-001
153	TR@FERN HILL	3/8/2016	7.90	7.20	-6.35E+	1.24	10.00	6.52	2.25	3.38	0.54	3.62	2.88	27.00	1.00E-001
154	TR@FERN HILL	3/22/2016	9.00	7.15	-6.35E+	1.33	10.00	7.30	2.47	3.85	0.72	4.11	3.19	29.20	1.00E-001
155	TR@FERN HILL	4/5/2016	10.50	7.19	-6.35E+	1.62	10.00	8.00	2.66	4.26	0.88	4.56	3.55	35.00	1.00E-001
156	TR@FERN HILL	4/13/2016	10.90	7.12	-6.35E+	1.40	10.00	8.32	2.75	4.45	0.96	4.76	3.85	34.20	1.00E-001
157	TR@FERN HILL	4/19/2016	15.20	7.21	-6.35E+	2.45	10.00	8.53	2.81	4.58	1.01	4.90	4.18	36.80	1.00E-001
158	TR@FERN HILL	4/26/2016	11.30	7.25	-6.35E+	1.40	10.00	8.84	2.89	4.76	1.08	5.09	4.42	36.10	1.00E-001
159	TR@FERN HILL	5/2/2016	14.30	7.23	-6.35E+	2.34	10.00	9.35	3.04	5.07	1.20	5.42	4.39	41.20	1.00E-001
160	TR@FERN HILL	5/16/2016	12.00	7.42	0.71	1.21	10.00	8.30	2.72	4.37	0.93	4.68	3.96	36.10	1.00E-001
161	TR@FERN HILL	6/13/2016	13.20	7.49	0.69	1.20	10.00	7.41	2.38	4.33	0.45	2.12	3.65	33.00	1.00E-001
162	TR@FERN HILL	6/27/2016	14.20	7.42	-6.35E+	1.12	10.00	6.98	2.38	3.65	0.65	3.90	3.52	29.70	1.00E-001
163	TR@FERN HILL	7/11/2016	12.30	7.57	0.66	1.19	10.00	6.16	2.15	3.66	0.37	2.12	3.57	29.20	1.00E-001
164	TR@FERN HILL	7/25/2016	11.80	7.27	-6.35E+	1.05	10.00	6.52	2.25	3.38	0.54	3.62	3.09	27.80	1.00E-001
165	TR@FERN HILL	8/15/2016	12.40	7.46	0.79	1.09	10.00	6.45	2.08	3.54	0.43	2.09	2.92	26.90	1.00E-001
166	TR@FERN HILL	8/29/2016	14.40	7.50	-6.35E+	1.12	10.00	6.48	2.24	3.36	0.53	3.59	3.00	27.10	1.00E-001
167	TR@FERN HILL	9/12/2016	15.70	7.50	-6.35E+	1.16	10.00	6.76	2.32	3.52	0.60	1.89	3.33	28.60	1.00E-001
168															





# Running the Model – Selecting Metal/Organism

BLM Freshwater version 3.1.2.37

File Edit Inputs Help

Description:

Current Selections

Prediction Mode: Toxicity Metal: C

Site Chemistry Simplified Site Chemistry

	Site Name	Sample Name	Temp.
			C



# Running the Model – Selecting Metal/Organism

BLM Freshwater version 3.1.2.37

File Edit Inputs Help

Description:

Current Selections

Prediction Mode: Toxicity Metal: C

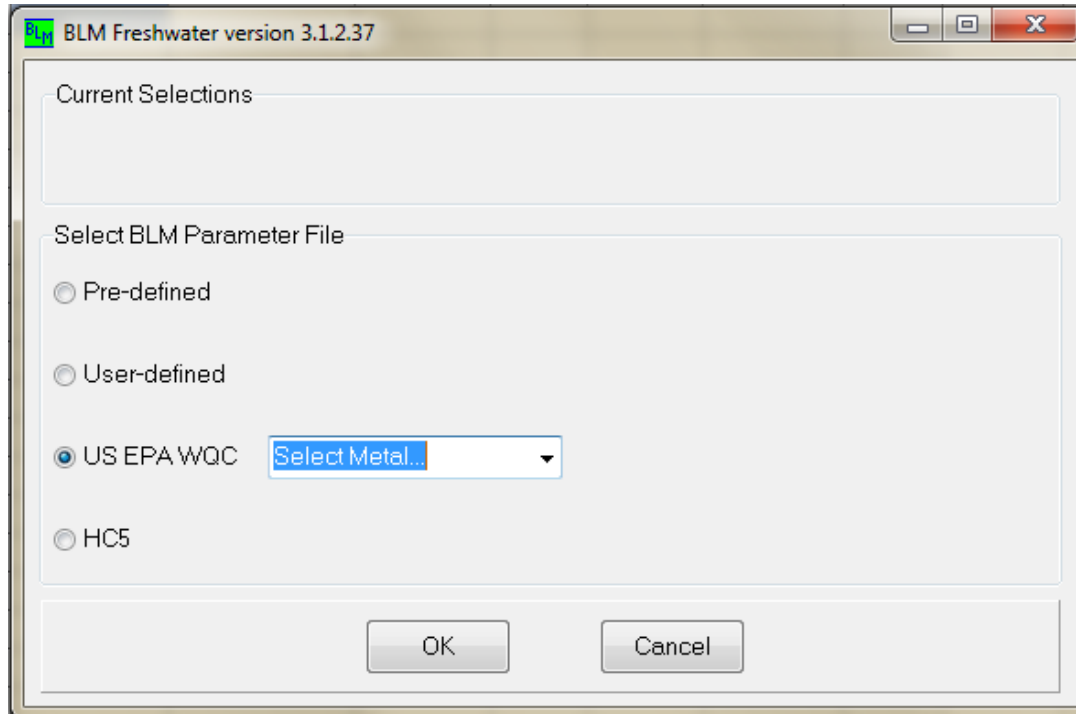
Site Chemistry Simplified Site Chemistry

	Site Name	Sample Name	Temp.
			C

Click to select metal/organism



# Selecting the Metal/Organism



BLM Freshwater version 3.1.2.37

Current Selections

Select BLM Parameter File

Pre-defined

User-defined

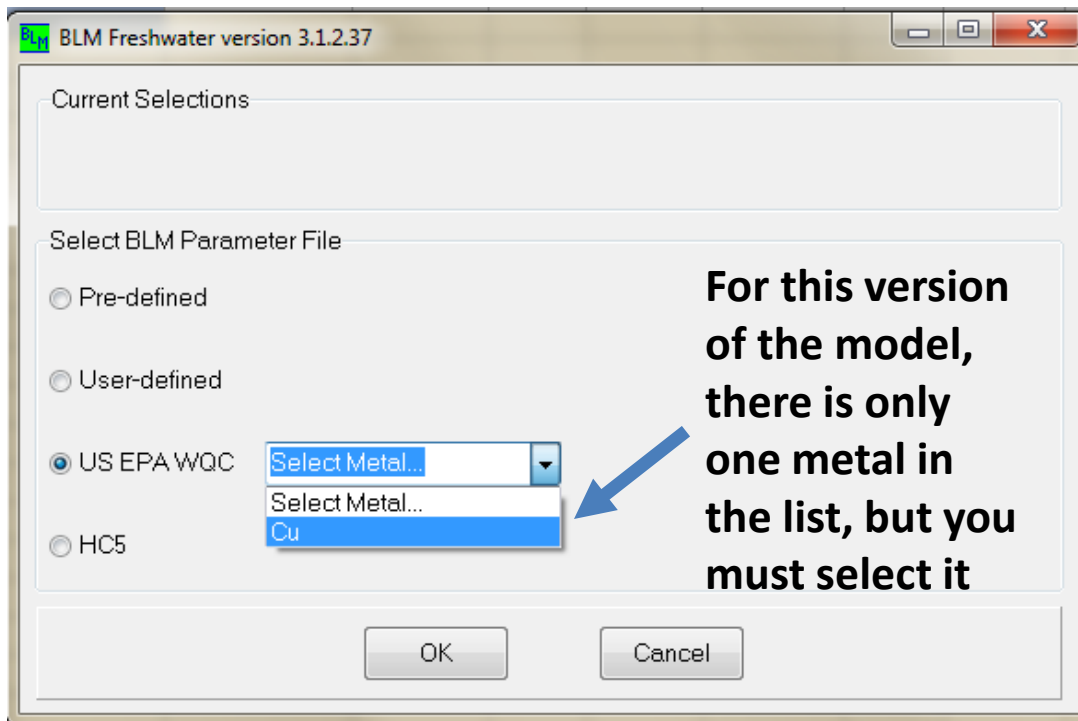
US EPA WQC

HC5

OK Cancel



# Selecting the Metal/Organism



# Running the Model

BLM Freshwater version 3.1.2.37

File Edit Inputs Help

Description:

Current Selections

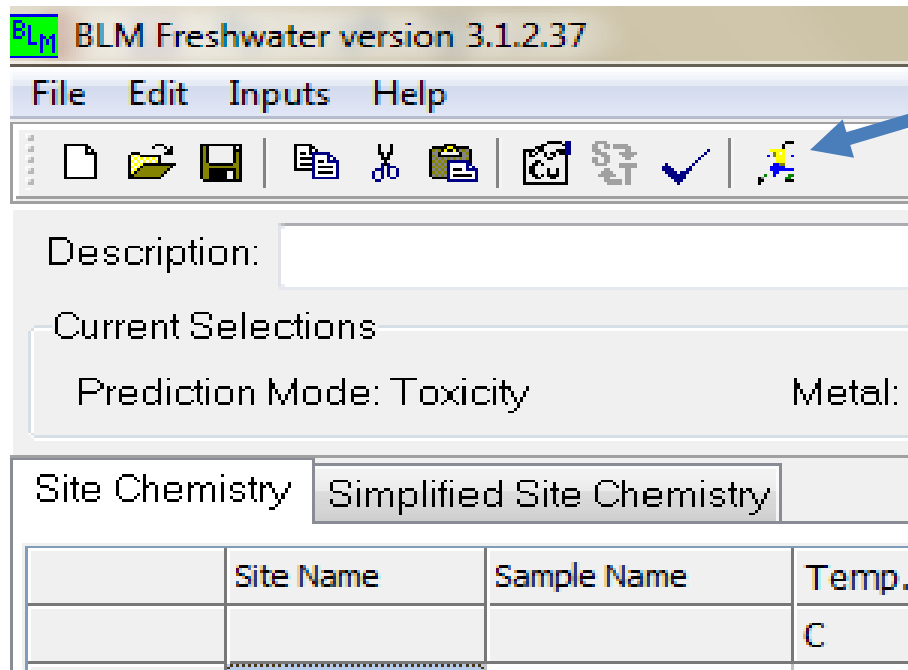
Prediction Mode: Toxicity Metal: C

Site Chemistry Simplified Site Chemistry

	Site Name	Sample Name	Temp.
			C



# Running the Model



The screenshot shows the BLM Freshwater software interface. The title bar reads "BLM Freshwater version 3.1.2.37". The menu bar includes "File", "Edit", "Inputs", and "Help". The toolbar contains icons for file operations (New, Open, Save, Print, Copy, Paste) and a "Run" button (a lightning bolt icon). A blue arrow points to the "Run" button. Below the toolbar, there is a "Description:" field, a "Current Selections" section with "Prediction Mode: Toxicity" and "Metal: C", and a "Site Chemistry" section with "Simplified Site Chemistry" selected. At the bottom, there is a table with columns for "Site Name", "Sample Name", and "Temp.".

	Site Name	Sample Name	Temp.
			C

**Click here to  
run the model**



# BLM Output – Water Quality Criteria

Site Label	Sample Label	Final Acute Value (FAV), ug/L	CMC (CMC=FAV/2), ug/L	CCC (CCC=FAV/ACR), ug/L	Cu ug/L	Acute Toxic Units (Acute TU=Cu/CMC)	Chronic Toxic Units (Chronic TU=Cu/CCC)	Censored Flag (0 = quantified, 1 = BDL)
"TR@FERN HILL	"1/19/2016	7.20	3.60	2.24	0.64	0.18	0.28	0.00
"TR@FERN HILL	"2/9/2016	5.24	2.62	1.63	0.62	0.24	0.38	0.00
"TR@FERN HILL	"2/23/2016	5.24	2.62	1.63	0.64	0.24	0.39	0.00
"TR@FERN HILL	"3/1/2016	4.42	2.21	1.37	0.79	0.36	0.58	0.00
"TR@FERN HILL	"3/8/2016	4.42	2.21	1.37	0.64	0.29	0.46	0.00
"TR@FERN HILL	"3/22/2016	4.42	2.21	1.37	0.64	0.29	0.46	0.00
"TR@FERN HILL	"4/5/2016	4.42	2.21	1.37	0.64	0.29	0.46	0.00
"TR@FERN HILL	"4/13/2016	4.42	2.21	1.37	0.64	0.29	0.46	0.00
"TR@FERN HILL	"4/19/2016	4.42	2.21	1.37	0.64	0.29	0.46	0.00
"TR@FERN HILL	"4/26/2016	4.42	2.21	1.37	0.64	0.29	0.46	0.00
"TR@FERN HILL	"5/2/2016	4.42	2.21	1.37	0.64	0.29	0.46	0.00
"TR@FERN HILL	"5/16/2016	7.80	3.90	2.42	0.71	0.18	0.30	0.00
"TR@FERN HILL	"6/13/2016	8.55	4.27	2.65	0.69	0.16	0.26	0.00
"TR@FERN HILL	"6/27/2016	8.55	4.27	2.65	0.64	0.15	0.24	0.00
"TR@FERN HILL	"7/11/2016	9.33	4.66	2.90	0.66	0.14	0.23	0.00
"TR@FERN HILL	"7/25/2016	9.33	4.66	2.90	0.64	0.14	0.22	0.00
"TR@FERN HILL	"8/15/2016	7.44	3.72	2.31	0.79	0.21	0.34	0.00
"TR@FERN HILL	"8/29/2016	7.44	3.72	2.31	0.64	0.17	0.27	0.00
"TR@FERN HILL	"9/12/2016	7.44	3.72	2.31	0.64	0.17	0.27	0.00



# Questions...

