## Data Submission Guide for CMDP Chemical/Radionuclide Samples (reported through EXCEL TEMPLATE)

reported through <u>LACLE TEMPLATE</u>

#### **Document Instructions**

The Excel template used for reporting Chemical/Radionuclide samples are separated into three sections including general *Sample Information* (see <u>page 2</u>), *Chem/Rad Results* (see <u>page 3</u>) and *Field Results and Measurements* (*Do not use for Chem/Rad reporting*). Additionally, this document includes the **Method Cart** (<u>pages 4-5</u>) and an **example** of a completed sample submission (<u>page 6</u>).

Submit the information as described on the following pages for **Chemical** and/or **Radionuclide** samples. This guidance document will highlight the required fields within each section.

Header fields in <u>RED and underlined text</u> are **required** in order to meet federal and state reporting requirements.
Header fields in <u>BLUE and underlined text</u> are *conditionally or situationally required*.
Header fields in **BLACK text** are not required for a successful data submission.
Sample data will be rejected if the <u>required fields</u> are left blank.

Additionally keep in mind:

- Use drop down icons ( ) to filter data.
- If you have questions please contact our CMDP support staff at <u>dec.cmdpsupport@alaska.gov</u>.

#### Screenshot of Chemical/Radionuclides Excel Template



## DATA SUBMISSION GUIDE FOR CHEMICAL/RADIONUCLIDE SAMPLES THROUGH EXCEL TEMPLATE

Header fields in **RED** and underlined text below are **required** in order to meet federal and state reporting requirements. Header fields in **BLUE** and underlined text are *conditionally or situationally required*. Header fields in **BLACK text** are not required for a successful data submission.

#### Section 1: Sample Information

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	C	MC	)P						Che	micals/F	Radion	uclides			
	Compliance	Monitori	ing Data Porta	al											
								I							
Reporti	ng Lab. ID *					Generate	XML								
							Sample I	nformatior	1 I						
Sample ID*	Sample Received Date <sup>f</sup>	WS ID*	Facility ID*	Sampling Point ID*	Sampling Location	Collection Date <sup>#<sup>f</sup></sup>	Collection Time (24H) <sup>f</sup>	Sample Type <sup>#f</sup>	Sample Volume	Repeat Location	Original Sample $ID^+$	Original Reporting Lab.ID	Original Collection Date	Comment	Sample Collector Name

**<u>Reporting Lab. ID</u>**: Lab identification number.

Sample ID: Lab sample identification number, limit to 20 characters (numbers, letters, dash/underscore (-\_) are allowed).

Sample Received Date: Date sample was received by lab (MM/DD/YY).

WS ID: Public Water System identification number (PWSID).

**NOTE:** Each water system name has a unique PWSID number. If the name and number on the work order does not match what is listed in CMDP, investigate to ensure accuracy and consistency.

Facility ID: Facility identification number where sample was collected.

<u>Sampling Point ID</u>: Sample Point identification number related to the facility where the sample was collected.

If you are unsure of the Facility ID and Sampling Point ID to enter, refer to either the:

- Monitoring Summary (PDF) which describes how to find this information through Drinking Water Watch.
- <u>PWS Facility/Sample Point List (XLS)</u> which provides the appropriate Facility ID and Sample Point ID for each analyte sampled for an individual public water system.

<u>Sampling Location</u>: This field must describe the location where the sample was taken (i.e., entry point, 123 Main Street, Health Clinic, etc.). Keep description succinct (*numbers, letters, dash/underscore* (-\_) *only*).

Collection Date: Date sample collected (MM/DD/YY).

Collection Time (24H): Time sample collected (HH:MM).

<u>Sample Type</u>: Type of sample to be submitted. From the drop down menu (pictured below), select the appropriate sample type (i.e., routine, confirmation, special, etc.).

Routine	
Repeat	-
Triggered	=
Confirmation	-
Special	
Batch Blanks	
Field Blanks	1
Performance Evaluation	*

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Sample Volume: Not required (numerical value only).

Repeat Location: Do NOT report data in this field.

<u>Original Sample ID</u>: If reporting a <u>confirmation</u> sample be sure to enter the lab <u>Sample ID</u> number of the original sample collected.

Original Reporting Lab. ID: Not required but report if information if available.

Original Collection Date: Not required but report if information if available (MM/DD/YY).

<u>Comment</u>: Must include water idle time when reporting Lead/Copper samples (e.g., Idle time 7 hours 40 minutes *or* Water sat unused for 9 hours). Please limit characters to numbers, letters, dash or underscore (-\_\_\_). In particular, do NOT include quotation marks.

Sample Collector Name: Name of sample collector, report if information is provided.

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Header fields in **RED** and underlined text below are required in order to meet federal and state reporting requirements. Header fields in **BLUE** and underlined text are conditionally or situationally required. Header fields in **BLACK text** are not required for a successful data submission.

#### Section 2: Chem/Rad Results

													_		
			_												
															( <i>I</i>
								Results							
Analyte <sup>#f</sup> [Code - Name]	Not Detected* <sup>f</sup>	Result <sup>f</sup>	Result UOM <sup>f</sup>	Standard Deviation (+/-) <sup>f</sup>	Reporting Limit <sup>f</sup>	Reporting Limit UOM	Volume Assayed (ML)	Method <sup>f</sup>	Analysis Start Date <sup>f</sup>	Analysis Start Time <sup>f</sup>	Analysis Completed Date	Analysis Completed Time	Analyst Name	Analyzing Lab ID	Comment
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Analyte [Code-Name]: From the drop down menu (pictured below), select the analytes you are submitting results for as appropriate.

0100-TURBIDITY	$\land$
0999-CHLORINE	
1000-TOTAL CHLORINE	
<b>1001-COMBINED CHLORIN</b>	
1002-ALUMINUM	
1003-NITROGEN-AMMONI	
1004-BROMIDE	
1005-ARSENIC	$\checkmark$

**NOTE:** This list will NOT filter by the analytes your lab (or lab that you have subcontracted with) is certified for so be sure to verify the certification information (outside of CMDP) prior to data submission.

Not Detected: From the drop down menu (pictured below), select whether the contaminant was Not Detected (select Yes) or **Detected** (select **No**) in the analyzed sample.

Yes No

**Result**: Enter the appropriate reported result of the sample only if reporting a detect.

Result UOM: From the drop down menu (pictured below), select the unit of measure for the sample result as appropriate.

MFL	^
pCi/L	
ng/L	
ug/L	
mg/L	
CU	
TON	
umho/cm	$\checkmark$

**Standard Deviation (+/-)**: Not required but report if applicable.

**Reporting Limit:** Enter the appropriate reporting limit of the analytes.

Reporting Limit UOM: From the drop down menu, select the unit of measure for the reporting limit as appropriate.

Volume Assayed: Not required but report if applicable (numerical value only).

Method: This drop down list includes methods for <u>ALL chemical/radionuclide analyses</u> and is not filtered by the analytes your lab (or subcontracting lab) is certified for. Be sure to verify certification status and select the appropriate method listed in the Methods Chart (see chart on pages 4-5). Notice that each Analysis Method Used by labs is assigned to a specific reporting code indicated in the Report this Method in CMDP column. Samples will be rejected if users report a code that is not listed on the methods chart.

Analysis Start Date: Date when lab began analysis (MM/DD/YY).

Analysis Start Time: Time when lab began analysis (HH:MM).

Analysis Completed Date: Not required but report if information is available (MM/DD/YY).

Analysis Completed Time: Not required but report if information is available (HH:MM).

Analyst Name: Not required.

Analyzing Lab ID: If the sample was subcontracted to a different lab for analysis, the analyzing lab identification number is required to be reported here.

Comment: Not required, however if comments are provided, please limit characters to numbers, letters, dash or underscore (- ). In particular, do NOT include quotation marks.

Section 3: Field Results and Measurements		Field Re	esults and	Measurements (Opt	ional)	
<b>REMINDER!</b> This section (pictured to right) should <b>NOT</b> be used for Chemical/Radionuclide sample submittals. Please leave fields blank	Parameter* [Code - Name]	Result*	Result UOM*	Method <sup>f</sup>	Analyst Name	Comment
chemical hadionuclide sample submittals. Please leave helds blank.						

## CHEMICAL/RADIONUCLIDE SAMPLES METHODS CHART

Analysis Method		
Used	Analyte	Report this Method in CDMP
	1,1,1-Trichloroethane-R, 1,1,2-Trichloroethane-R, 1,1-	
	Dichloroethylene-R, 1,2,4-Trichlorobenzene-R, 1,2-Dichlorobenzene-R,	
	1,2-Dichloroethane-K, 1,2-Dichloropropane-K, 1,4-Dichlorobenzene-K, Benzene-R. Bromodichloromethane-T. Bromoform-T. Carbon	
524.2	Tetrachloride-R. Chlorobenzene-R. Chloroform-T. cis-1.2-	524.2-VOC. GC/MS. P&T. CAPCOLUMN
52 1.2	Dichloroethylene-R. Dibromochloromethane-T. Dichloromethane	
	(Methylene Chloride)-R, Ethylbenzene-R, Styrene-R,	
	Tetrachloroethylene-R, Toluene-R, Total THM-T, Total Xylenes-R,	
	trans-1,2 Dichloroethylene, Trichloroethylene-R, Vinyl Chloride-R	
504.1	1,2-Dibromo-3-chloropropane-DBCP, 1,2-Dibromoethane-EDB	504.1-GC-MICROEXTRACTION-ECD
515.3	2,4,5-TP, 2,4-D, Dalapon, Dicamba, Dinoseb, Pentachlorophenol (PCP),	
	B-Hydroxycarbofuran Aldicarb Aldicarb Sulfone Aldicarb Sulfoxide	515.5-515.5
531.2	Carbaryl, Carbofuran, Methomyl, Oxamyl (vydate)	531.2-531.2
	Acenapthene, Acenapthylene, Alachlor, Aldrin, Atrazine,	
	Benzo(a)pyrene, Butachlor, Di(2-ethylhexyl)adipate, Di(2-	
525.2	ethylhexyl)phthalate, Dieldrin, Endrin, Heptachlor, Heptachlor	
525.2	epoxide, Hexachlorobenzene, Hexachlorocyclopentadiene (HCCPD),	
	Lindane (g-BHC), Methoxychlor, Metolachlor, Metribuzin, Propachlor,	
2220D (20th Ed.)	Simazine	
2320B (2011 EQ.)	Alkalinity	2320B-TITRIMETRIC
2320B (21st Ed.)	Alkalinity	2320B-TITRIMETRIC
2320B-97 (online)	Alkalinity	2320B-TITRIMETRIC
	Aluminum, Antimony, Arsenic, Barium, Beryllium, Cadmium,	
200.8	Chromium, Copper, Lead, Manganese, Nickel, Selenium, Silver,	
	Thallium, Uranium, Zinc	200.8-INDUCTIVELY COUPLED PLASMA MASS SPECTROM 200.8
100.2	Asbestos	100.2-TRANSMISSION ELECTRON MICROSCOPY
300.1	Bromate	300.1-ION CHROMATOGRAPHY
317	Bromate, Chlorite	317-317
300.0	Bromide, Chlorate, Chloride, Chlorite, Fluoride, Nitrate-N, Nitrite-N,	
	Calcium Iron Magnesium Potassium Sodium Aluminum Barium	
200.7	Bervllium, Cadmium, Calcium, Chromium, Copper, Iron, Magnesium,	
	Manganese, Nickel, Potassium, Silver, Sodium	200.7-INDUCTIVELY COUPLED PLASMA
505	Chlordane, PCB monitoring, Toxaphene	505-PESTICIDES, PCB, GC, MICROEXTRACT
2120B (20th Ed.)	Color	2120B-VISUAL COMPARISON METHOD
2120B (21st Ed.)	Color	2120B-VISUAL COMPARISON METHOD
2120B (22nd Ed.)	Color	2120B-VISUAL COMPARISON METHOD
2510B (20th Ed.)	Conductivity	2510B-CONDUCTANCE @ 25C
2510B (21st Ed.)	Conductivity	2510B-CONDUCTANCE @ 25C
2510B (22110 EU.)	Conductivity	
2310B-97 (Online)	Conductivity	2310B-CONDUCTANCE @ 25C
2330B (20° Ed.)	Corrosivity	2330B - 2330B 2330B - 2330B
335.4	Cvanide	335 4-SPECTROPHOTOMETRIC SEMI-AUTOMATED
4500-CN-C (21st Ed.)	Cyanide	4500CN-C-AMENABLE SPECTROPHOTOMETRIC
4500-CN-C (online)	Cyanide	4500CN-C-AMENABLE SPECTROPHOTOMETRIC
4500-CN-E (21st Ed.)	Cyanide	4500CN-E-SPECTROPHOTOMETRIC, MANUAL
4500-CN-E (online)	Cyanide	4500CN-E-SPECTROPHOTOMETRIC, MANUAL
4500-CN-C, E (online)	Cyanide	4500CN-E - SPECTROPHOTOMETRIC, MANUAL
552.2	Dibromoacetic acid, Dichloroacetic acid, Monobromoacetic acid,	552.2-DBPS & CL2 SOLVENTS GC L/L ELECTRON CAPT
	Monochloroacetic acid, Total Haloacetic Acids, Trichloroacetic acid	
552.3	Ulbromoacetic acid, Dichloroacetic acid, Monopromoacetic acid,	
1613	Dioxin (2.3.7.8-TCDD)	1613-HIGH RES GCMS-CAPCOL-1 /L EXTRACTION
549.2	Diquat. Paraguat	549.2-549.2
5310B (21st Ed.)	Dissolved Organic Carbon (DOC), Total Organic Carbon (TOC)	5310B-HIGH TEMPERATURE COMBUSTION METHOD
5310C (20th Ed.)	Dissolved Organic Carbon (DOC), Total Organic Carbon (TOC)	5310C-PERSULFATE-ULTRAVIOLET OR OXIDATION
5310C (21st Ed.)	Dissolved Organic Carbon (DOC), Total Organic Carbon (TOC)	5310C-PERSULFATE-ULTRAVIOLET OR OXIDATION
5310C (22 <sup>nd</sup> Ed.)	Dissolved Organic Carbon (DOC), Total Organic Carbon (TOC)	5310B-HIGH TEMPERATURE COMBUSTION METHOD
548.1	Endothall	548.1-GC-L/S EXTRACTION-ELECTRON CAPTURE DETEC
4500-F-C (22nd Ed.)	Fluoride	4500F-C-POTENTIOMETER ION SELECTIVE ELECTRODE
4500-F-C (online)	Fluoride	4500F-C-POTENTIOMETER ION SELECTIVE ELECTRODE
901.1	Gamma/Photon emitters	
547	Giypnosphate	p47-HIGH PERF LIQ CHROIVI-POST COL REACTOR-FLU

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## CHEMICAL/RADIONUCLIDE SAMPLES METHODS CHART

Analysis Method		
Used	Analyte	Report this Method in CDMP
7110B (22nd Ed.)	Gross alpha evaporation, Gross beta evaporation	7110B-EVAPORATION METHOD FOR GROSS ALPHA-BETA
900.0	Gross alpha, Gross beta	900 -EVAPORATION METHOD FOR GROSS ALPHA-BETA
2340B (20th Ed.)	Hardness	2340B-HARDNESS IN WATER BY CALCULATION
2340B (22nd Ed.)	Hardness	2340B-HARDNESS IN WATER BY CALCULATION
2340B (online.)	Hardness	2340B-HARDNESS IN WATER BY CALCULATION
5540C (20th Ed.)	MBAS-Foaming Agents	5540C-Methylene Blue Active Substances (MBAS)
245.1	Mercury	245.1-MANUAL COLD VAPOR TECHNIQUE
4500-NO3-D (20th Ed.)	Nitrate	4500NO3-D-ION SELECTIVE ELECTRODE
4500-NO3-E (20th Ed.)	Nitrate-N, Nitrate-Nitrite as N	4500NO3-E-CADMIUM REDUCTION, MANUAL
HACH 10206 (Jan. 2011)	Nitrate-N, Nitrate-Nitrite as N	HACH 10206-HACH 10206
353.2	Nitrate-N, Nitrate-Nitrite as N, Nitrite-N	353.2-CADMIUM REDUCTION, AUTOMATED
4500-NO3-F (21st Ed.)	Nitrate-N, Nitrite-N	4500NO3-F-CADMIUM REDUCTION, AUTOMATED
4500-NO2-B (20th Ed.)	Nitrite-N	4500NO2-B-SPECTROPHOTOMETRIC, MANUAL
4500-NO2-B (online)	Nitrite-N	4500NO2-B - SPECTROPHOTOMETRIC, MANUAL
2150B (20 <sup>th</sup> Ed.)	Odor	2150B - Threshold Odor Test
2150B (21st Ed.)	Odor	2150B-Threshold Odor Test
4500-P-E (20th Ed.)	Ortho-phosphate	4500P-E-COLORIMETRIC, MANUAL
4500-P-E (21st Ed.)	Ortho-phosphate	4500P-E-COLORIMETRIC, MANUAL
4500-P-E (22 <sup>nd</sup> Ed.)	Ortho-phosphate	4500P-E-COLORIMETRIC, MANUAL
150.1	рН	150.1-ELECTROMETRIC-INDIVIDUAL MEASUREMENT
4500-H-B (20th Ed.)	рН	4500H-B-ELECTROMETRIC-ONLINE MEASUREMENT
4500-H-B (21st Ed.)	рН	4500H-B-ELECTROMETRIC-ONLINE MEASUREMENT
4500-H-B-00 (online)	рН	4500H-B-ELECTROMETRIC-ONLINE MEASUREMENT
7500Ra-B (22nd Ed.)	Radium 226	7500-RAB-PRECIPITATION METHOD FOR RADIUM
7500Ra-D (22nd Ed.)	Radium 228	7500-RAD-SEQUENTIAL PRECIPITATION METHOD FOR RADI
903.0	Radium-226	903.0-PRECIPITATION METHOD FOR RADIUM
903.1	Radium-226	903.1-PRECIPITATION METHOD FOR RADIUM
904.0	Radium-228	904.0-SEQUENTIAL PRECIPITATION METHOD FOR RADI
905.0	Strontium-89, 90	905.0-LIQUID SCINTILLATION SPECTROPHOTOMETRIC
2540C (20th Ed.)	TDS	2540C-Total Dissolved Solids Dried at 180 deg C
2540C (21st Ed.)	TDS	2540C-Total Dissolved Solids Dried at 180 deg C
2540C (22nd Ed.)	TDS	2540C-Total Dissolved Solids Dried at 180 deg C
2540C-97 (online)	TDS	2540C-Total Dissolved Solids Dried at 180 deg C
906.0	Tritium LSC	906-LIQUID SCINTILLATION SPECTROPHOTOMETRIC
180.1	Turbidity	180.1-180.1
2130B (20th Ed.)	Turbidity	2130B-NEPHELOMETRIC
2130B (21st Ed.)	Turbidity	2130B-NEPHELOMETRIC
5910B (20 <sup>th</sup> Ed.)	UV 254	5910B-ULTRAVIOLET ABSORPTION METHOD
5910B (22nd Ed.)	UV 254	5910B-ULTRAVIOLET ABSORPTION METHOD

# DATA SUBMISSION GUIDE FOR CHEMICAL/RADIONUCLIDE SAMPLES THROUGH EXCEL TEMPLATE

### Section 4: Example of Completed Sample Submission

1	Lab. ID *	AK009	0961			G	Generate XI	ИL								6																			
<b>–</b>						Sample	Informatio	n																Results								F	ield Results an	d Measurements (Opt	ional)
Sample ID*	Sample Received Date <sup>f</sup>	WS ID*	Facility <sup>!</sup> ID*	Sampling Point ID*	mpling C	Collection Date <sup>*f</sup>	Collection Time (24H) <sup>f</sup>	Sample Type <sup>*f</sup>	Sample Volume (ML)	Repeat Location	Original Sample ID <sup>+</sup>	Original Reporting Lab.ID	Original Collection Date	Comment	Sample Collector Name	Analyte* <sup>f</sup> [Code - Name]	Not Detected *	Result	f Result	Standard Deviation (+/-) <sup>f</sup>	Reporting Limit <sup>f</sup>	Reporting Limit UOM <sup>+/</sup>	g Volume Assayed (ML)	Method <sup>f</sup>	Analysis Start Date <sup>f</sup>	Analysis Start Time <sup>f</sup>	Analysis Completed Date	Analysis Completed Time	Analyst Name	Analyzing Lab ID	Comment	Parame Code - N	DO	NOT USE	omme
80-1612	2 5/10/2020	AK2263080	30 TP001	SPTP001 Ent	ry Point (	05/10/20	10:10	Routine							КВ	2378-1,2,4-TRICHLOROBENZENE	Yes				0.5	5 ug/L		524.2-VOC, GC/MS, P&T, CAPCOLUMN	5/11/2020	08:00							FOR	CHEN/RΔΓ	
																380-CIS-1,2-DICHLOROETHYLENE	Yes				0.5	5 ug/L		524.2-VOC, GC/MS, P&T, CAPCOLUMN	5/11/2020	08:00							rony		
																955-XYLENES, TOTAL	No	4.6	ug/L		0.9	5 ug/L		524.2-VOC, GC/MS, P&T, CAPCOLUMN	5/11/2020	08:00						<b></b> •			

# **Sample Information**

1 ortin	g Lab. ID *	AK009	61			G	ienerate XI	ИL							
						Sample	Informatio	n							
Sample ID*	Sample Received Date <sup>f</sup>	WS ID*	Facility ID*	Sampling Point ID*	Sampling Location	Collection Date <sup>*<sup>f</sup></sup>	Collection Time (24H) <sup>f</sup>	Sample Type <sup>*f</sup>	Sample Volume (ML)	Repeat Location	Original Sample ID <sup>+</sup>	Original Reporting Lab.ID	Original Collection Date	Comment	Sample Collector Name
280-1612	5/10/2020	AK2263080	TP001	SPTP001	Entry Point	05/10/20	10:10	Routine							KB

# **Example of a Sample Result (Non-Detect):**

						· · · · · · · · · · · · · · · · · · ·			/				1	
							Results							
Analyte <sup>#f</sup> [Code - Name]	Not Detected* <sup>f</sup>	Result <sup>f</sup> Result UOM <sup>f</sup>	Standard Deviation (+/-) <sup>f</sup>	Reporting Limit <sup>f</sup>	Reporting Limit UOM <sup>+f</sup>	Volume Assayed (ML)	Method <sup>f</sup>	Analysis Start Date <sup>f</sup>	Analysis Start Time <sup>f</sup>	Analysis Completed Date	Analysis Completed Time	Analyst Name	Analyzing Lab ID	Comment
2378-1,2,4-TRICHLOROBENZENE	Yes	<u> </u>	//	0.5	ug/L	1/	524.2-VOC, GC/MS, P&T, CAPCOLUMN	5/11/2020	08:00	<u> </u>	1	1		1
2380-CIS-1,2-DICHLOROETHYLENE	Yes	· · · · · · · · · · · · · · · · · · ·	,	0.5	ug/L	4'	524.2-VOC, GC/MS, P&T, CAPCOLUMN	5/11/2020	08:00					1
,				,		,		1		1	1	1	1	1

# **Example of a Sample Result (Detect):**

$\overline{\mathcal{O}}$															
								Results							
Analyte <sup>#f</sup> [Code - Name]	Not Detected* <sup>f</sup>	Result <sup>f</sup>	Result UOM <sup>f</sup>	Standard Deviation (+/-) <sup>f</sup>	Reporting Limit <sup>f</sup>	Reporting Limit UOM <sup>+f</sup>	Volume Assayed (ML)	Method <sup>f</sup>	Analysis Start Date <sup>f</sup>	Analysis Start Time <sup>f</sup>	Analysis Completed Date	Analysis Completed Time	Analyst Name	Analyzing Lab ID	Comment
2955-XYLENES, TOTAL	No	4.6	ug/L		0.5	ug/L		524.2-VOC, GC/MS, P&T, CAPCOLUMN	5/11/2020	08:00					