

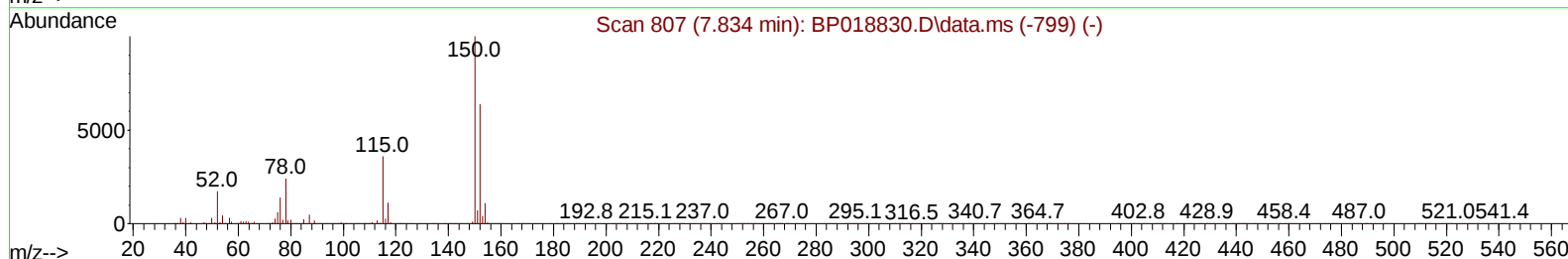
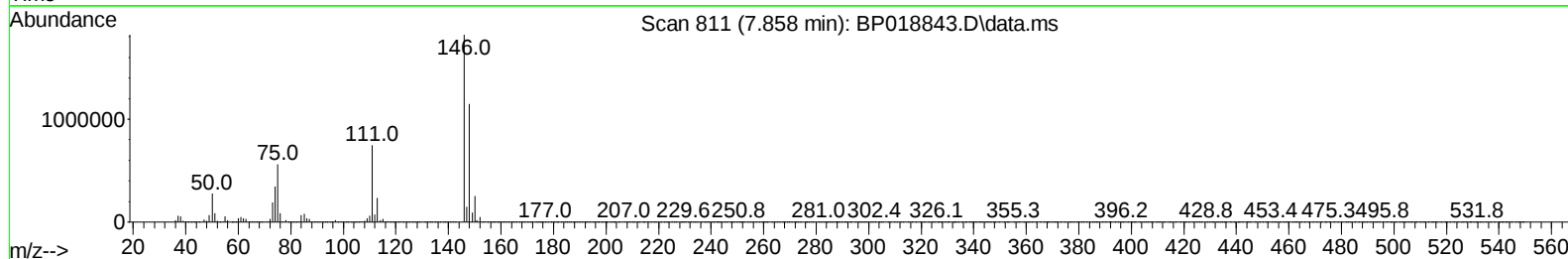
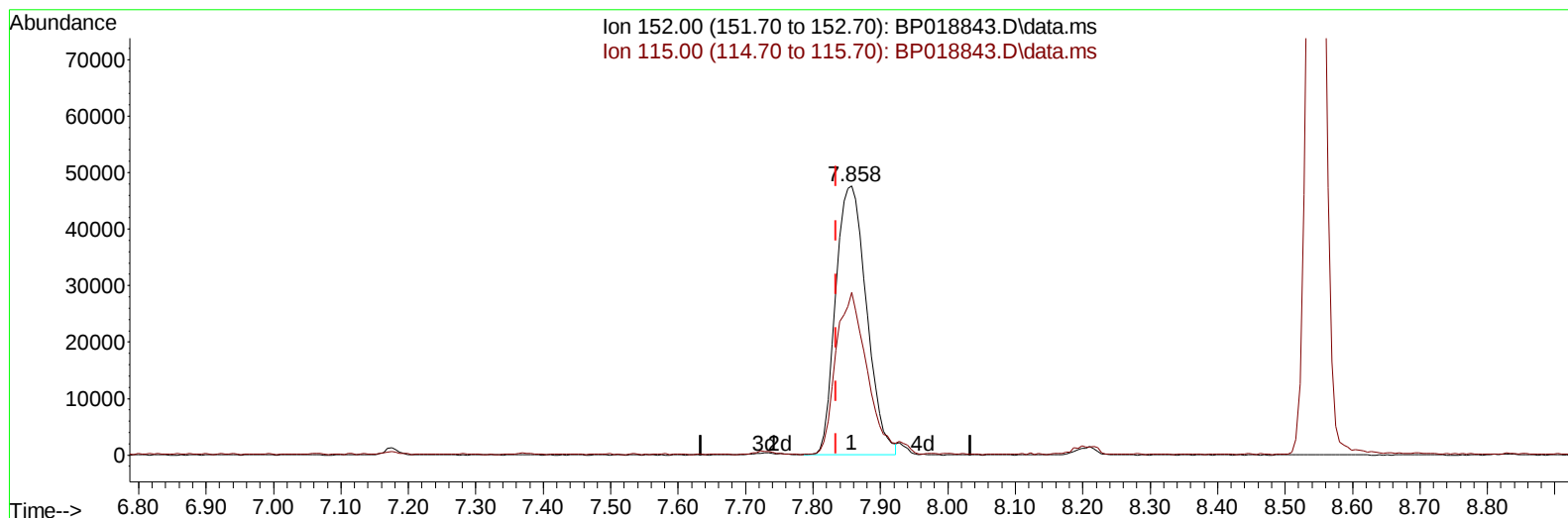
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP121323\
 Data File : BP018843.D
 Acq On : 14 Dec 2023 15:19
 Operator : MA/JU
 Sample : 05646-16
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AS3

Manual IntegrationsAPPROVED

Quant Time: Dec 14 21:54:17 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Dec 13 23:34:47 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/15/2023
 Supervised By :mohammad ahmed 12/16/2023



TIC: BP018843.D\data.ms

(1) 1,4-Dichlorobenzene-d4 (I)

7.858min (+ 0.024) 20.00 ng/u1

response 151224

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	59.50	60.49
0.00	0.00	0.00
0.00	0.00	0.00

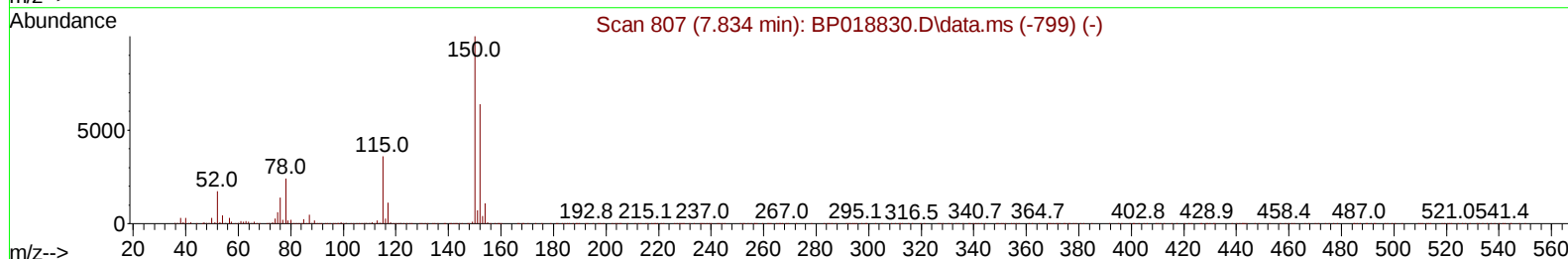
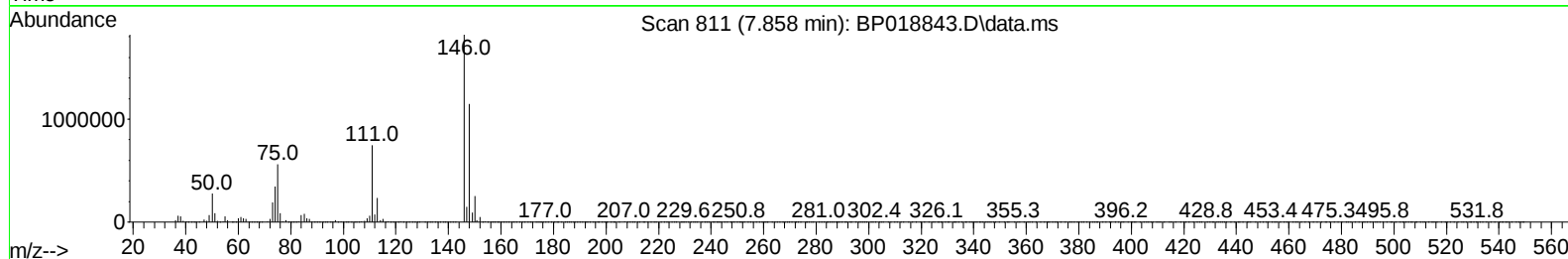
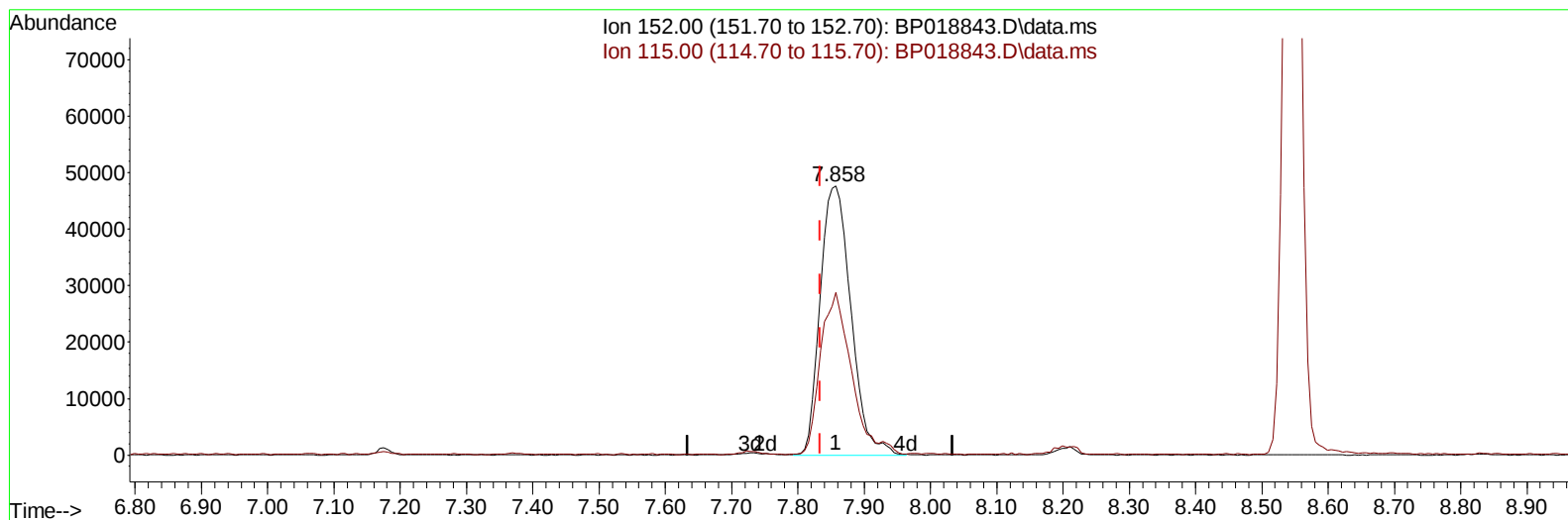
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TIC: BP018843.D\data.ms

(1) 1,4-Dichlorobenzene-d4 (I)

7.858min (+ 0.024) 20.00 ng/ul m

response 153919

Ion	Exp%	Act%
152.00	100.00	100.00
115.00	59.50	60.49
0.00	0.00	0.00
0.00	0.00	0.00

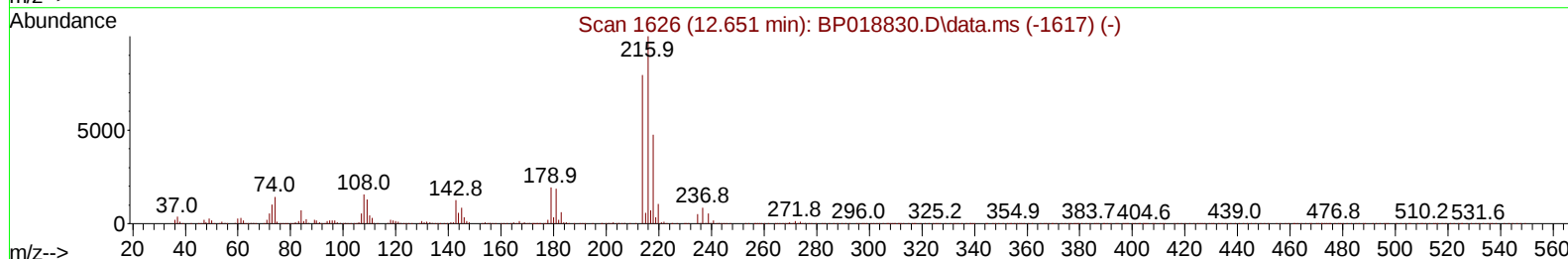
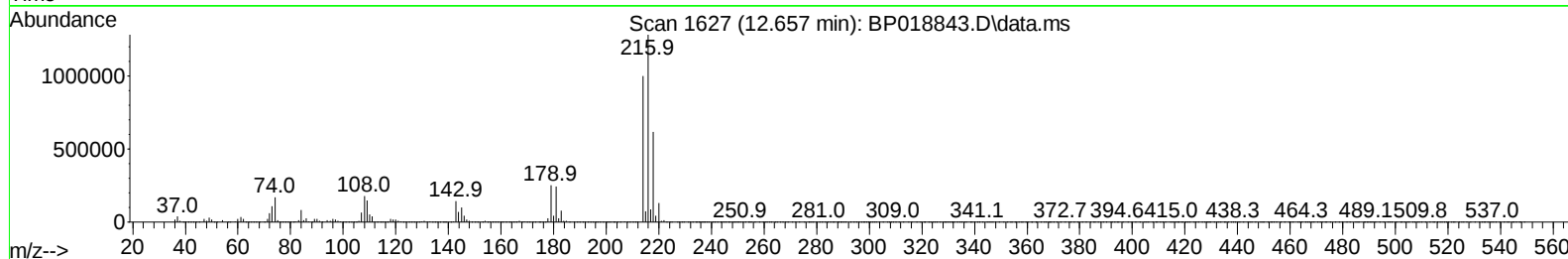
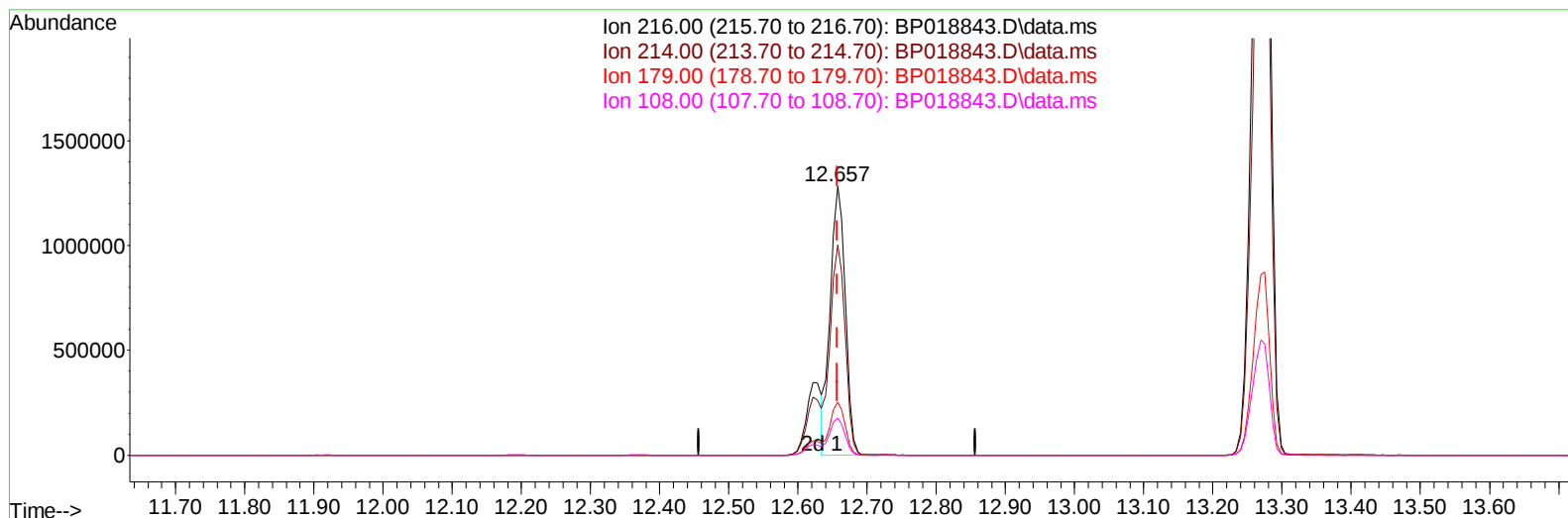
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 Sample : 05646-16
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Manual IntegrationsAPPROVED

Quant Time: Dec 14 22:24:34 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
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TIC: BP018843.D\data.ms

(39) 1,2,4,5-Tetrachlorobenzene

12.657min (+ 0.000) 147.53 ng/u1

response 1940229

Ion	Exp%	Act%
216.00	100.00	100.00
214.00	79.90	78.14
179.00	20.40	19.76
108.00	15.80	13.89

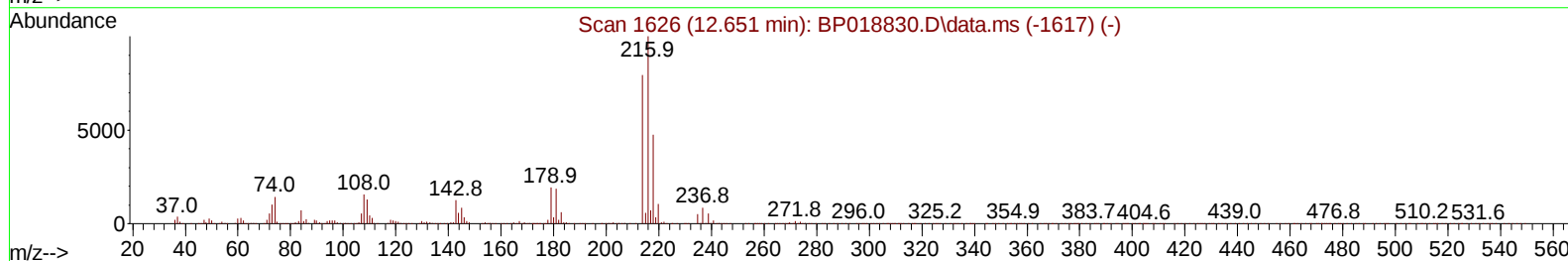
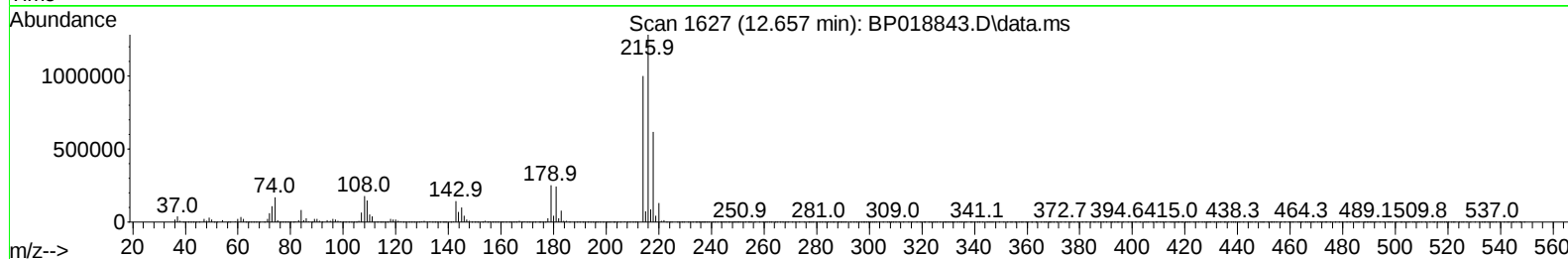
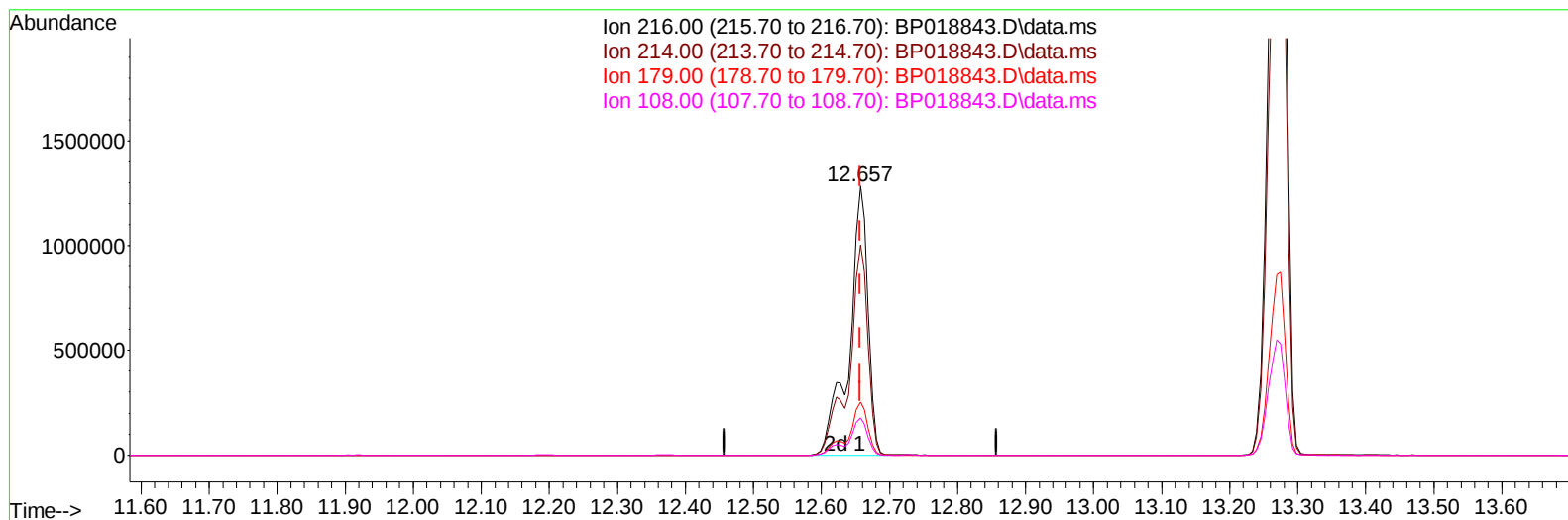
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TIC: BP018843.D\data.ms

(39) 1,2,4,5-Tetrachlorobenzene

12.657min (+ 0.000) 187.79 ng/ul m

response 2469766

Ion	Exp%	Act%
216.00	100.00	100.00
214.00	79.90	78.14
179.00	20.40	19.76
108.00	15.80	13.89

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/15/2023
 Supervised By :mohammad ahmed 12/16/2023

Quant Time: Dec 14 22:25:24 2023
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 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	153919m	20.000	ng/ul	0.02
20) Naphthalene-d8	10.657	136	607393	20.000	ng/ul	0.03
38) Acenaphthene-d10	14.463	164	362494	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.210	188	821596	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	967509	20.000	ng/ul	0.00
88) Perylene-d12	23.657	264	1176986	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	26890	5.953	ng/uL	0.00
4) Pyridine-d5	3.681	84	315114	26.046	ng/ul	0.00
7) Phenol-d5	7.005	99	464408	30.066	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.175	67	293127	32.019	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	386065	32.031	ng/ul	0.00
15) 4-Methylphenol-d8	8.546	113	352832	28.226	ng/ul	0.00
21) Nitrobenzene-d5	8.999	128	171955	31.808	ng/ul	0.00
24) 2-Nitrophenol-d4	9.716	143	182796	32.614	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.263	165	367489	32.066	ng/ul	0.01
31) 4-Chloroaniline-d4	10.781	131	455189	30.391	ng/ul	0.01
46) Dimethylphthalate-d6	13.881	166	933869	28.865	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	1133958	31.830	ng/ul	0.00
54) 4-Nitrophenol-d4	14.669	143	127060	23.926	ng/ul	0.00
60) Fluorene-d10	15.451	176	876869	32.289	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	98595	20.014	ng/ul	0.00
73) Anthracene-d10	17.310	188	1379147	31.930	ng/ul	0.00
81) Pyrene-d10	19.545	212	1810583	24.041	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.504	264	2145926	31.270	ng/ul	-0.01
Target Compounds						
					Qvalue	
8) Phenol	7.034	94	130585	8.621	ng/ul	100
18) 4-Methylphenol	8.605	108	198682	15.690	ng/ul	87
39) 1,2,4,5-Tetrachloroben...	12.657	216	2469766m	187.789	ng/ul	
69) Hexachlorobenzene	16.510	284	15970	1.379	ng/ul	98
75) 1,2,3,4-Tetrachloroben...	13.275	216	6987217	491.844	ng/uL	98
76) Pentachlorobenzene	14.781	250	782382	56.066	ng/uL	99

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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