

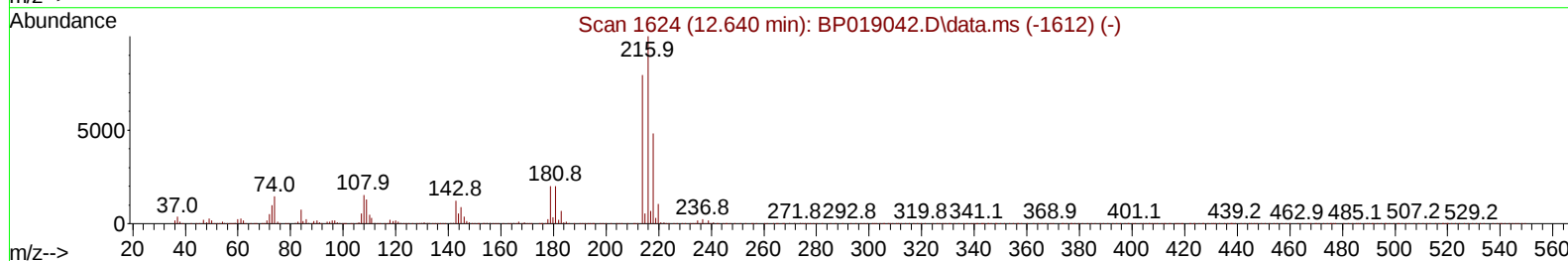
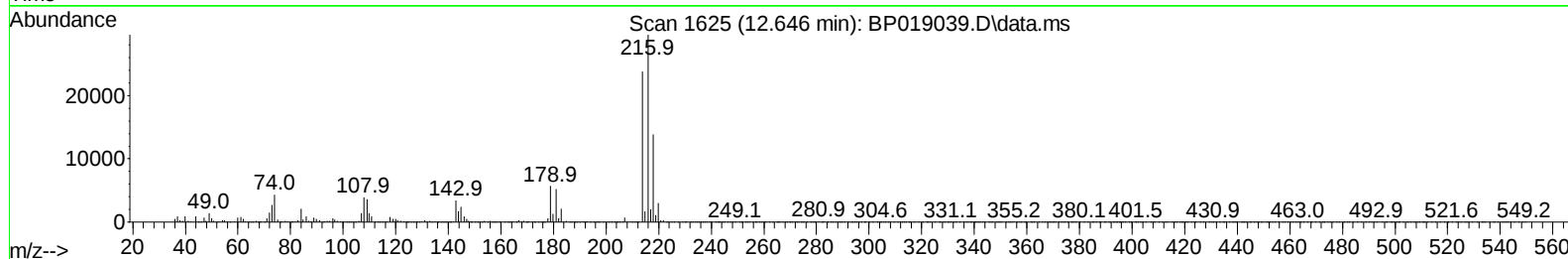
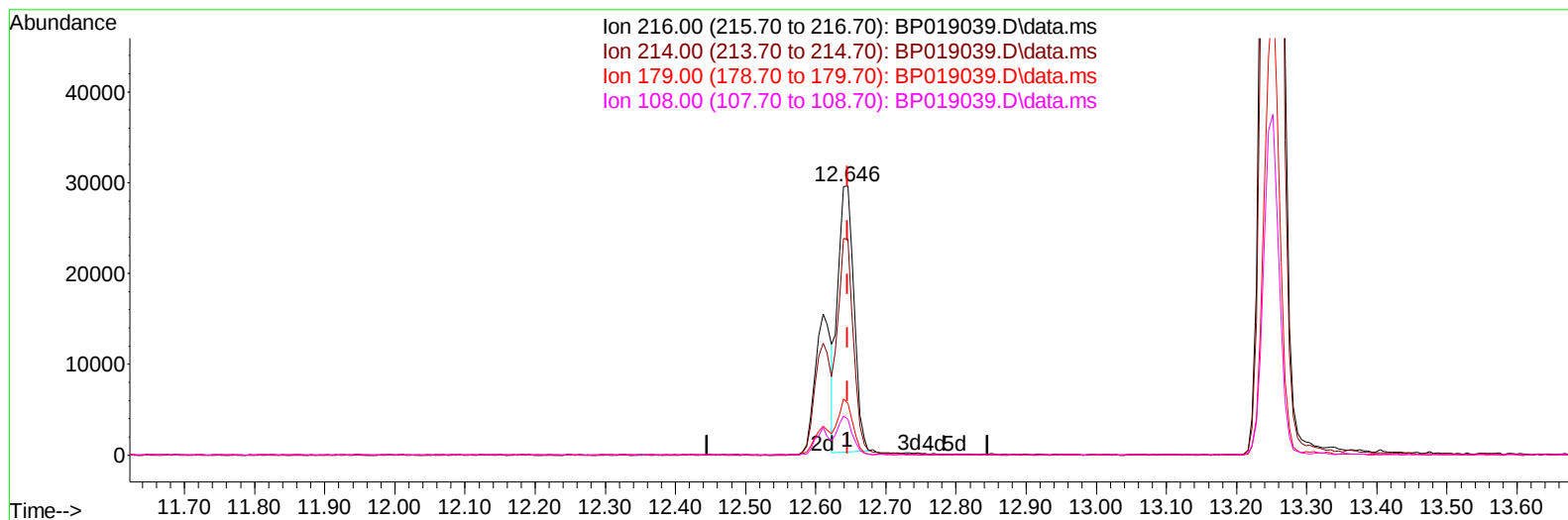
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122123\
 Data File : BP019039.D
 Acq On : 22 Dec 2023 14:18
 Operator : MA/JU
 Sample : 05808-09DL 5X
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B03DL

Manual IntegrationsAPPROVED

Quant Time: Dec 22 14:45:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 21 04:46:12 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 12/25/2023
 Supervised By :mohammad ahmed 12/25/2023



TIC: BP019039.D\data.ms

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

12.646min (+ 0.000) 3.44 ng/u1

response 46877

Ion	Exp%	Act%
216.00	100.00	100.00
214.00	79.90	80.37
179.00	20.40	19.17
108.00	15.80	13.29

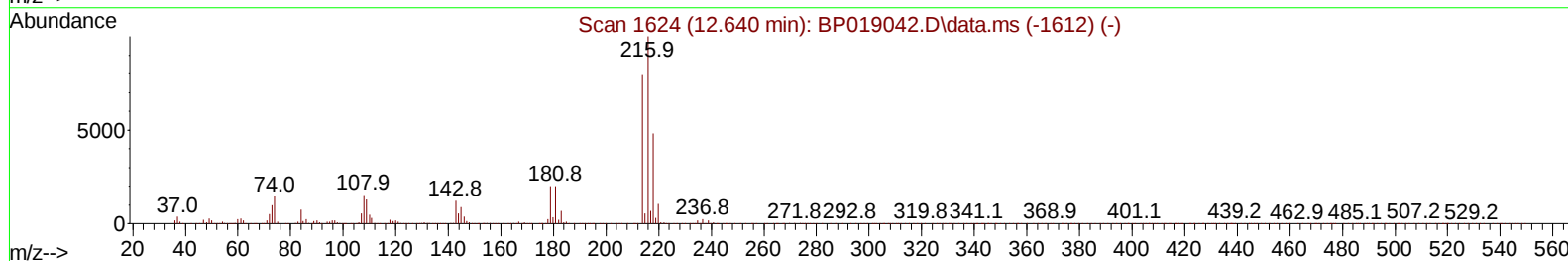
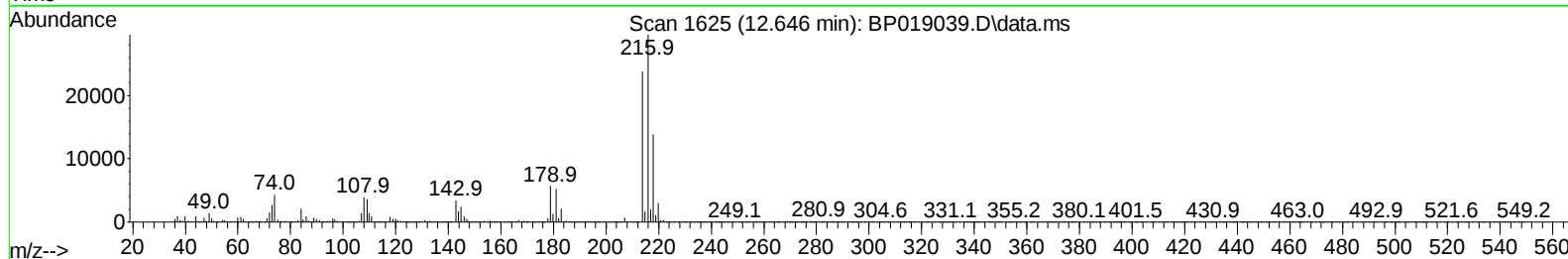
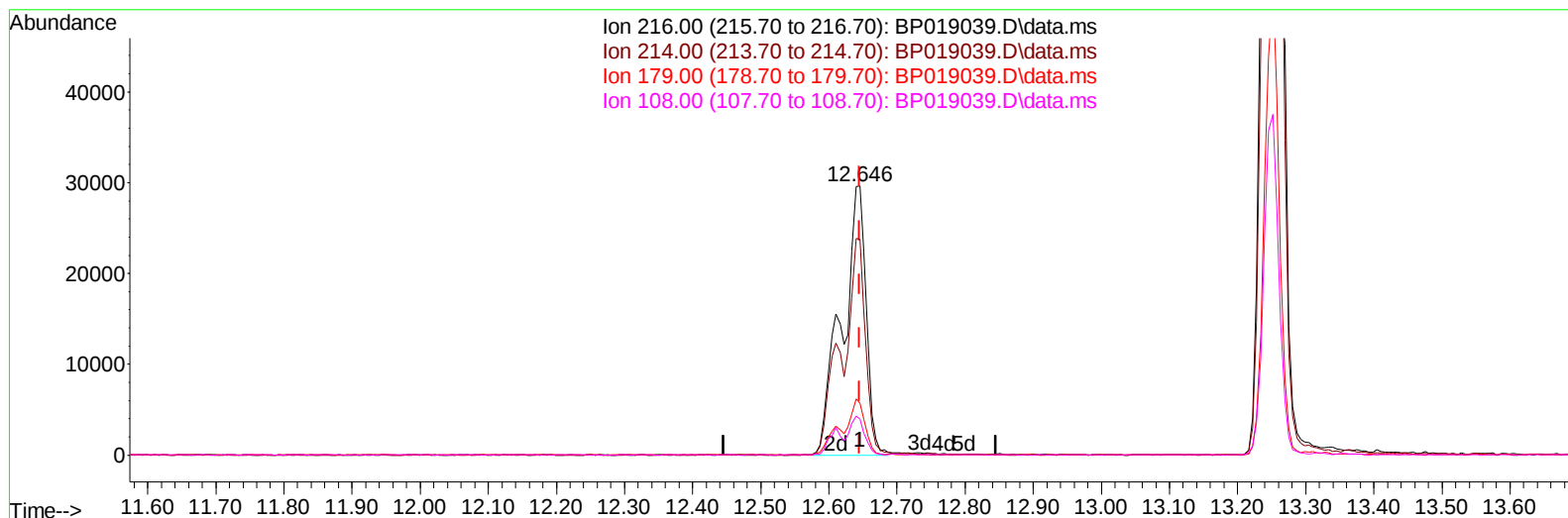
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TIC: BP019039.D\data.ms

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

12.646min (+ 0.000) 5.35 ng/ul m

response 72876

Ion	Exp%	Act%
216.00	100.00	100.00
214.00	79.90	80.37
179.00	20.40	19.17
108.00	15.80	13.29

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 12/25/2023
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Quant Time: Dec 22 22:53:06 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120123.MA.M
 Quant Title : SVOA CALIBRATION
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	153997	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	573899	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.451	164	375547	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	899777	20.000	ng/ul	0.00
79) Chrysene-d12	21.298	240	974472	20.000	ng/ul	0.00
88) Perylene-d12	23.657	264	1137871	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	3827	0.847	ng/uL	0.00
4) Pyridine-d5	3.681	84	29995	2.478	ng/ul	0.01
7) Phenol-d5	6.993	99	66271	4.288	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.158	67	40196	4.388	ng/ul	0.00
11) 2-Chlorophenol-d4	7.352	132	51483	4.269	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	50010	3.999	ng/ul	0.00
21) Nitrobenzene-d5	8.981	128	23485	4.598	ng/ul	0.00
24) 2-Nitrophenol-d4	9.699	143	24469	4.621	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	48082	4.440	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	60985	4.309	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	163229	4.870	ng/ul	0.00
49) Acenaphthylene-d8	14.146	160	177567	4.811	ng/ul	0.00
54) 4-Nitrophenol-d4	14.687	143	17311	3.146	ng/ul	0.02
60) Fluorene-d10	15.445	176	152287	5.413	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.575	200	17035	3.158	ng/ul	0.00
73) Anthracene-d10	17.304	188	246295	5.207	ng/ul	0.00
81) Pyrene-d10	19.545	212	311817	4.111	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.504	264	327812	4.941	ng/ul	0.00
Target Compounds						
					Qvalue	
39) 1,2,4,5-Tetrachloroben...	12.646	216	72876m	5.349	ng/ul	
75) 1,2,3,4-Tetrachloroben...	13.251	216	381406	24.515	ng/uL	98
76) Pentachlorobenzene	14.769	250	72627	4.752	ng/uL	100

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

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