

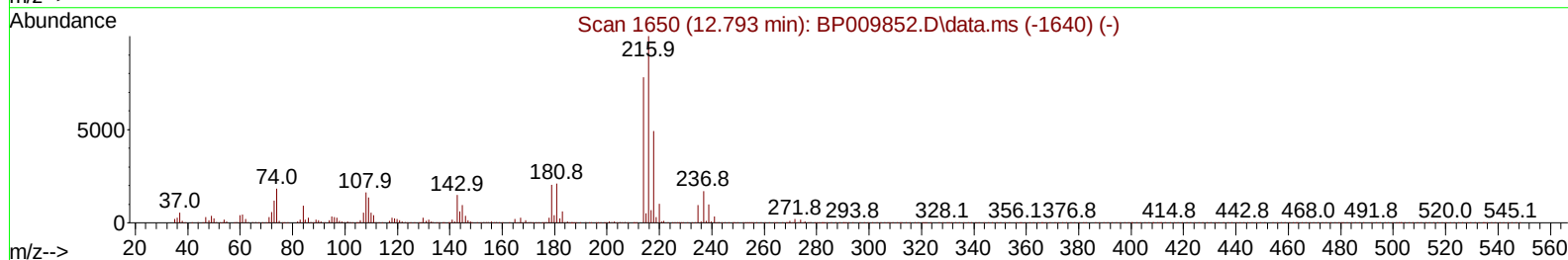
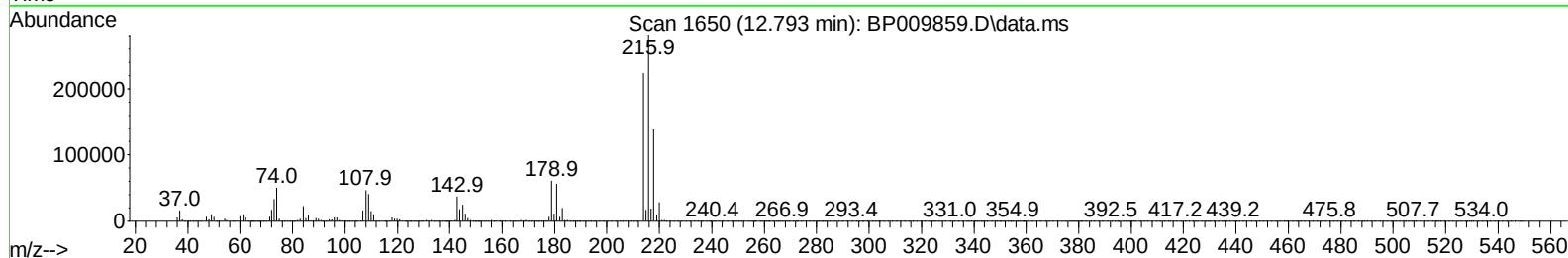
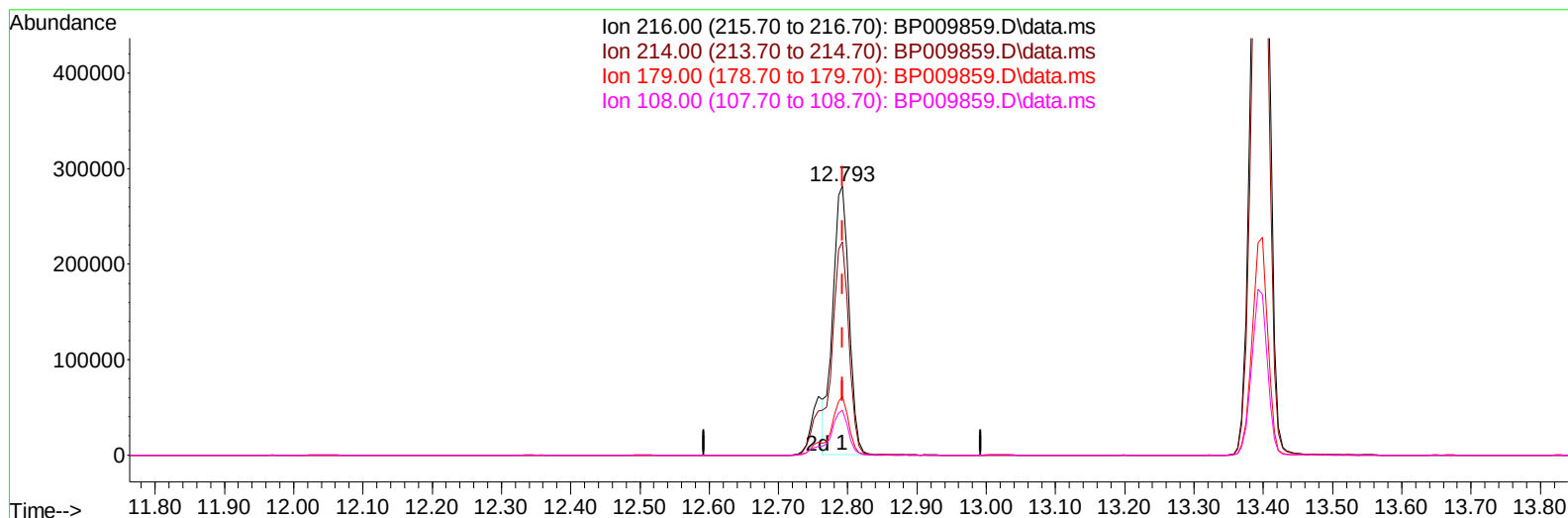
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP041522\
 Data File : BP009859.D
 Acq On : 15 Apr 2022 13:56
 Operator : CG/JU
 Sample : N2422-05
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0J47

Manual IntegrationsAPPROVED

Quant Time: Apr 18 01:01:01 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP040722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Apr 18 00:59:30 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/18/2022
 Supervised By :Yogesh Patel 04/18/2022



TIC: BP009859.D\data.ms

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

12.793min (+ 0.000) 36.51 ng/ul

response 455713

Ion	Exp%	Act%
216.00	100.00	100.00
214.00	72.70	79.38
179.00	17.70	21.74#
108.00	19.10	16.72

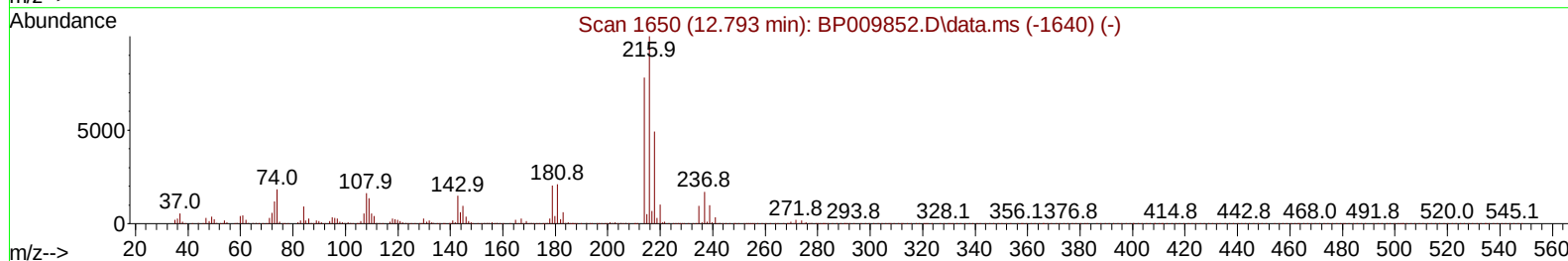
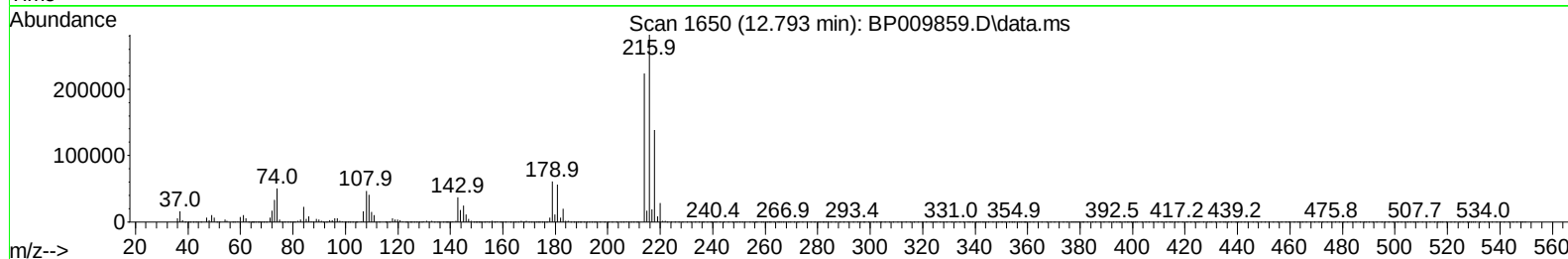
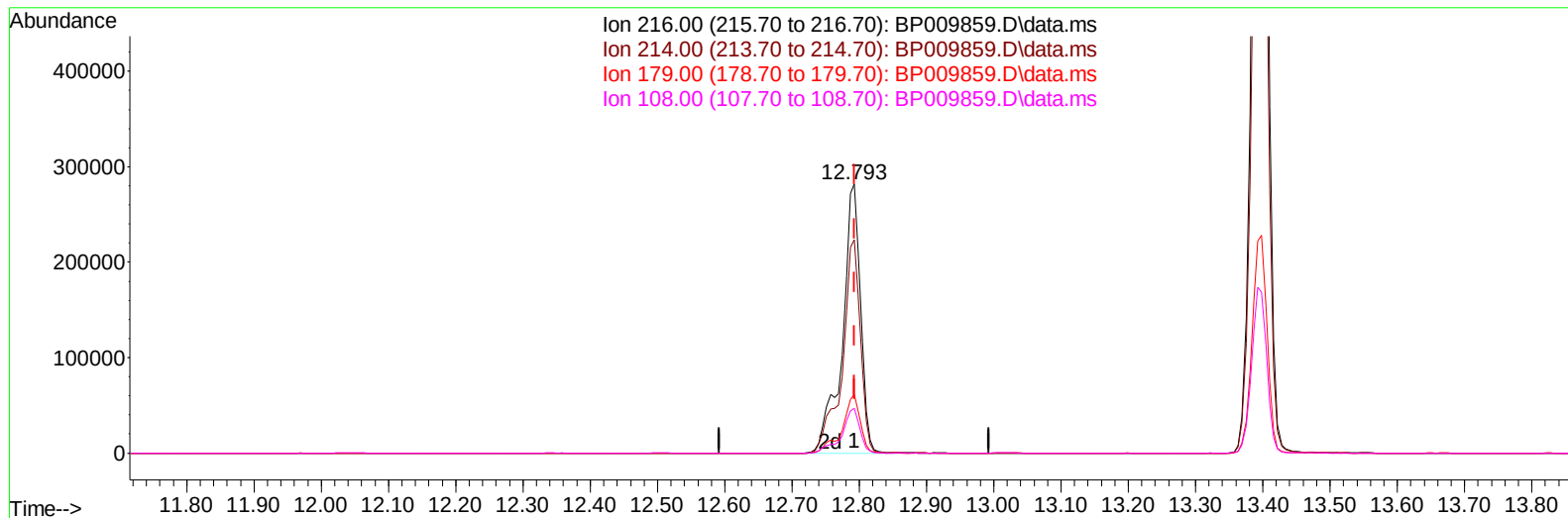
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TIC: BP009859.D\data.ms

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

12.793min (+ 0.000) 42.65 ng/ul m

response 532344

Ion	Exp%	Act%
216.00	100.00	100.00
214.00	72.70	79.38
179.00	17.70	21.74#
108.00	19.10	16.72

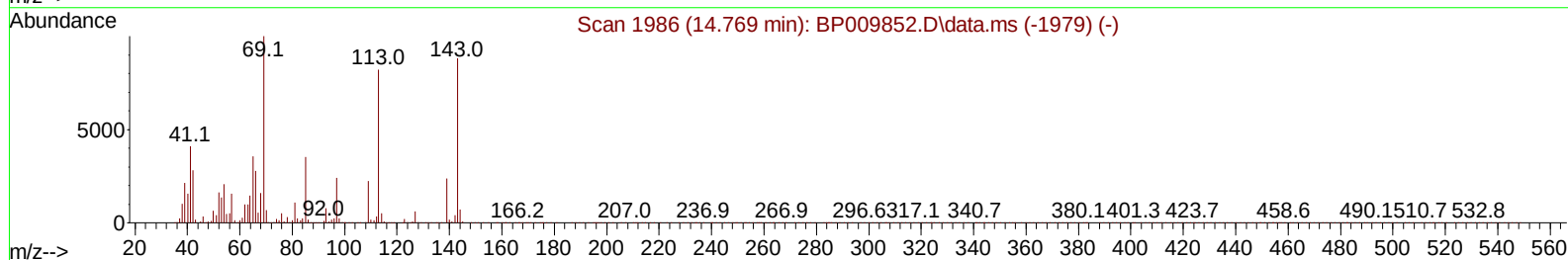
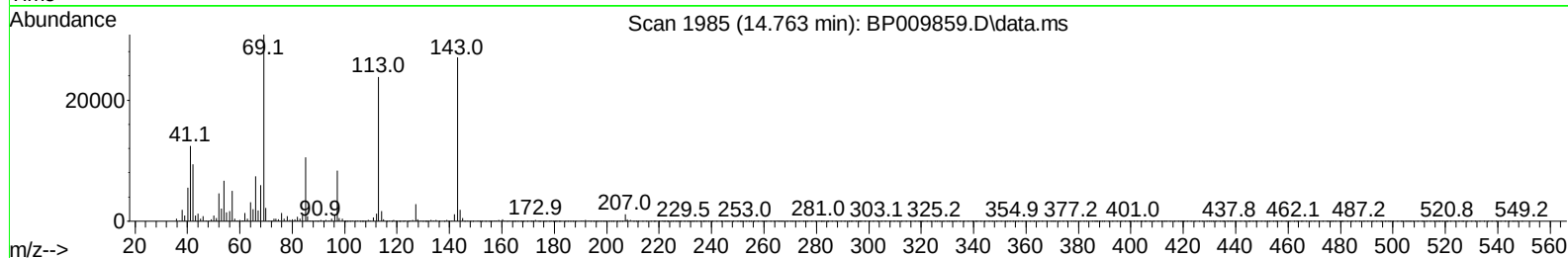
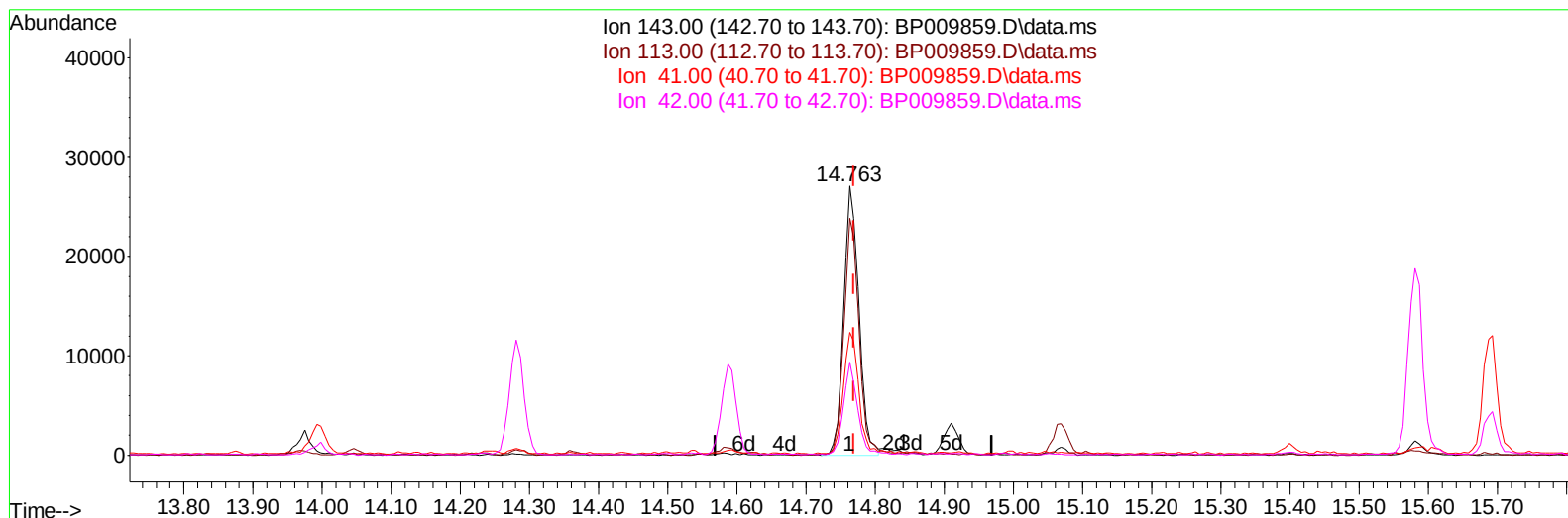
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TIC: BP009859.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.006) 7.16 ng/u1

response 41593

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	88.10#
41.00	23.20	45.72#
42.00	18.00	34.62#

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.975	152	135336	20.000	ng/ul	# 0.02
20) Naphthalene-d8	10.787	136	542073	20.000	ng/ul	0.02
38) Acenaphthene-d10	14.592	164	407981	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.339	188	944827	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.422	240	1002805	20.000	ng/ul	# 0.00
88) Perylene-d12	23.886	264	935851	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	13133	4.331	ng/uL	0.00
4) Pyridine-d5	3.805	84	68114	8.147	ng/ul	0.00
7) Phenol-d5	7.116	99	78718	6.625	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.287	67	237896	33.632	ng/ul	0.00
11) 2-Chlorophenol-d4	7.493	132	227368	25.369	ng/ul	0.00
15) 4-Methylphenol-d8	8.663	113	156392	15.767	ng/ul	0.00
21) Nitrobenzene-d5	9.122	128	156525	40.042	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143	167575	39.101	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.387	165	278371	31.040	ng/ul	0.00
31) 4-Chloroaniline-d4	10.904	131	231535	18.127	ng/ul	0.01
46) Dimethylphthalate-d6	13.998	166	1150330	35.895	ng/ul	0.00
49) Acenaphthylene-d8	14.281	160	1304168	34.350	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	42741m	7.354	ng/ul	0.00
60) Fluorene-d10	15.581	176	1006585	37.337	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.692	200	189836	33.916	ng/ul	0.00
73) Anthracene-d10	17.439	188	1683648	37.218	ng/ul	0.00
81) Pyrene-d10	19.669	212	2131343	38.916	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.727	264	1934798	39.700	ng/ul	0.00
Target Compounds						
						Qvalue
22) Nitrobenzene	9.163	77	743342	74.085	ng/ul#	85
39) 1,2,4,5-Tetrachloroben...	12.793	216	532344m	42.648	ng/ul	
75) 1,2,3,4-Tetrachloroben...	13.398	216	1615162	118.239	ng/uL#	86
76) Pentachlorobenzene	14.910	250	38404	2.884	ng/uL#	86

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

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