

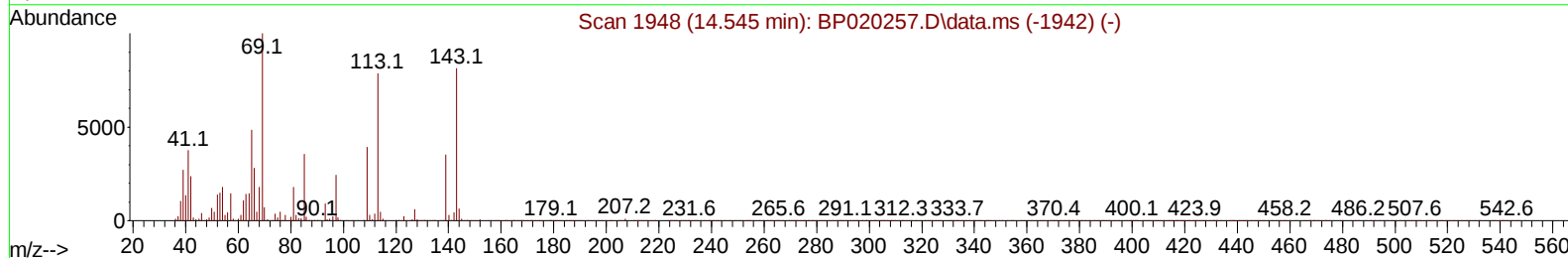
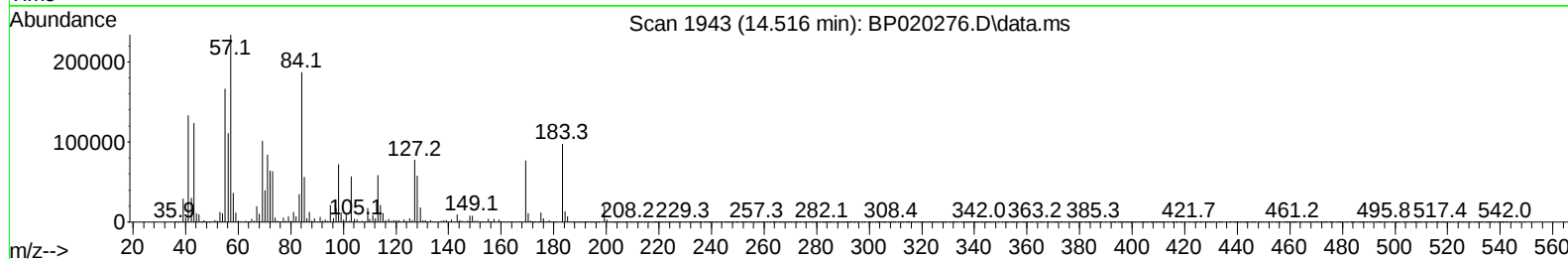
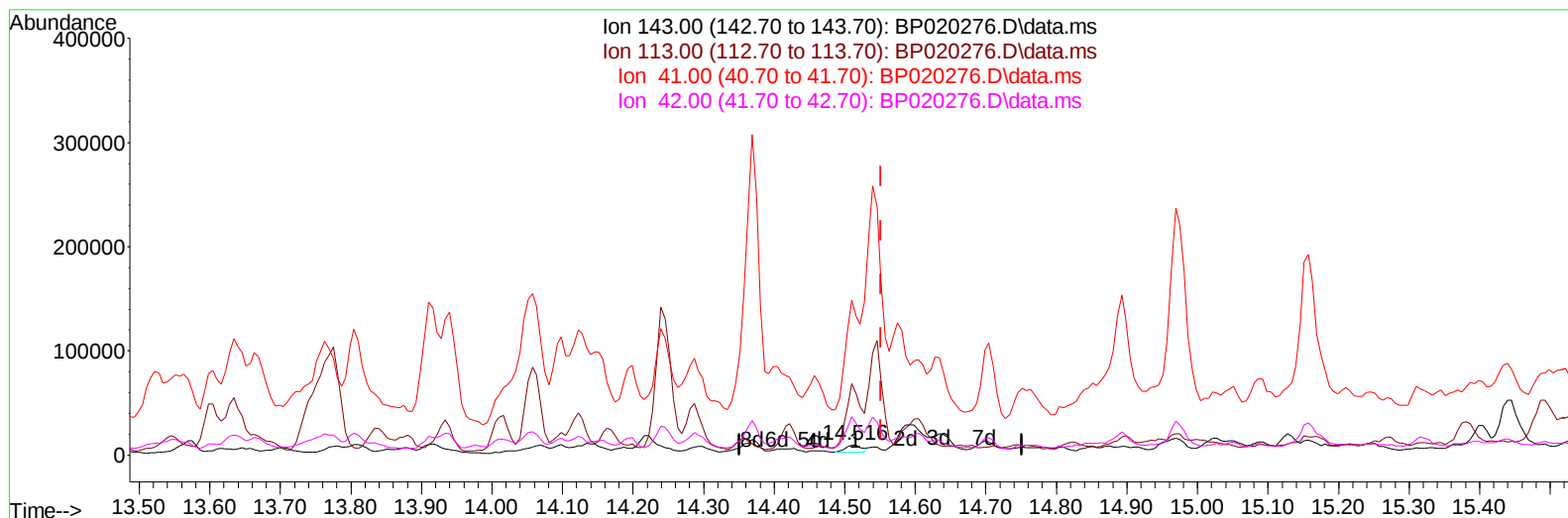
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050824\
 Data File : BP020276.D
 Acq On : 09 May 2024 21:10
 Operator : MA/JU
 Sample : P2369-12
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A43N6

Manual IntegrationsAPPROVED

Quant Time: May 09 22:45:47 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed May 08 22:57:55 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/10/2024
 Supervised By :mohammad ahmed 05/11/2024



TIC: BP020276.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (-0.035) 3.90 ng/u1

response 10938

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	101.40	607.35#
41.00	50.30	1369.12#
42.00	32.50	313.94#

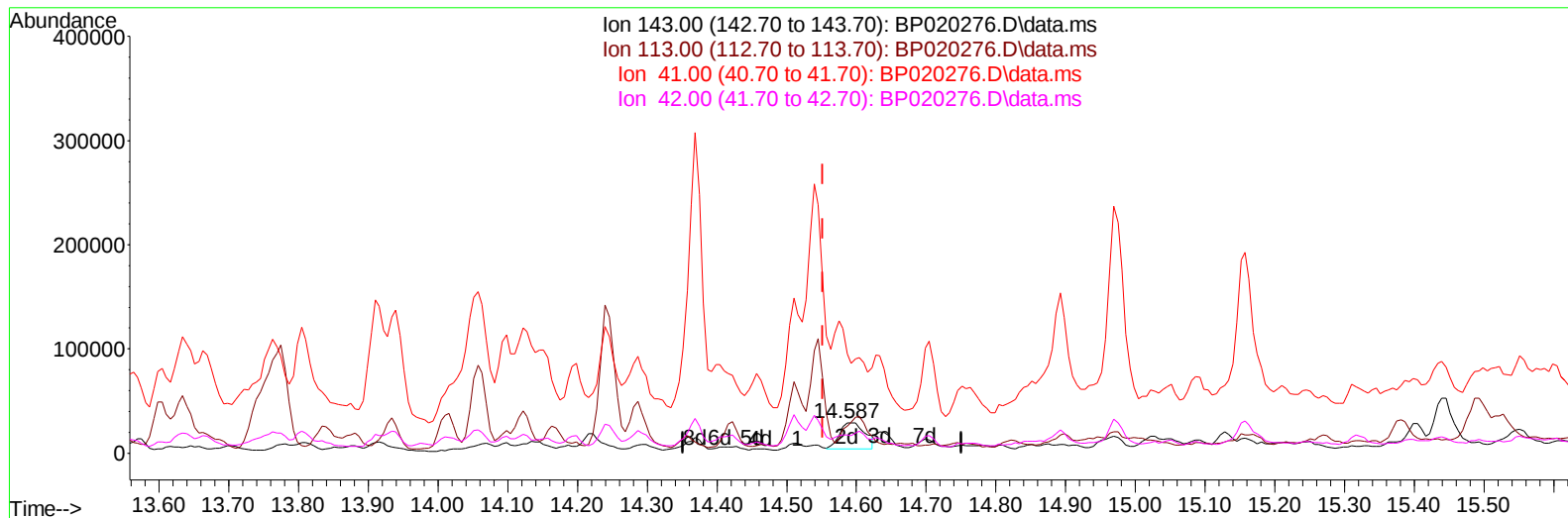
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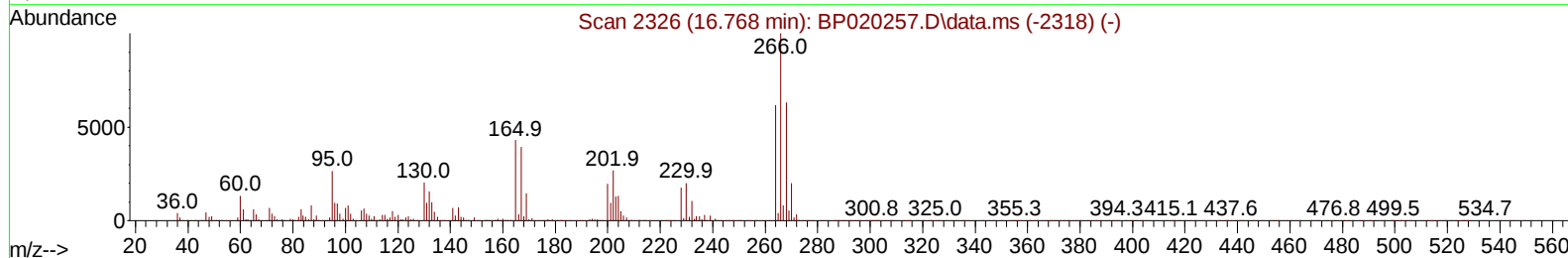
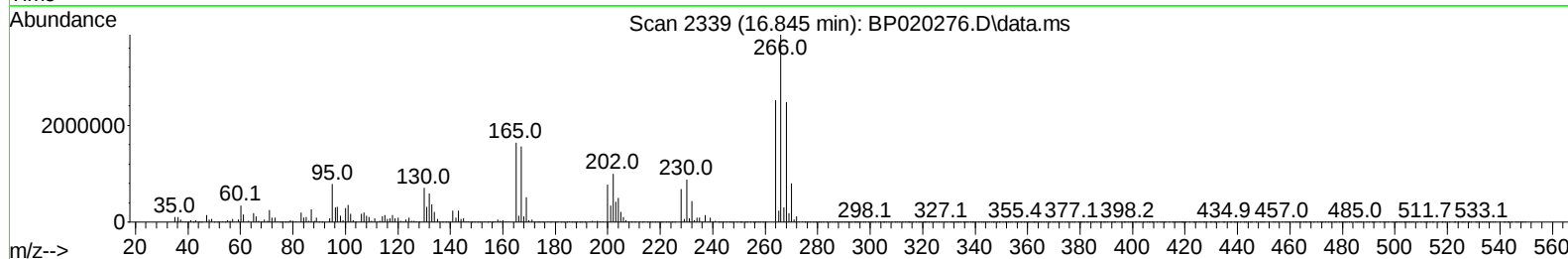
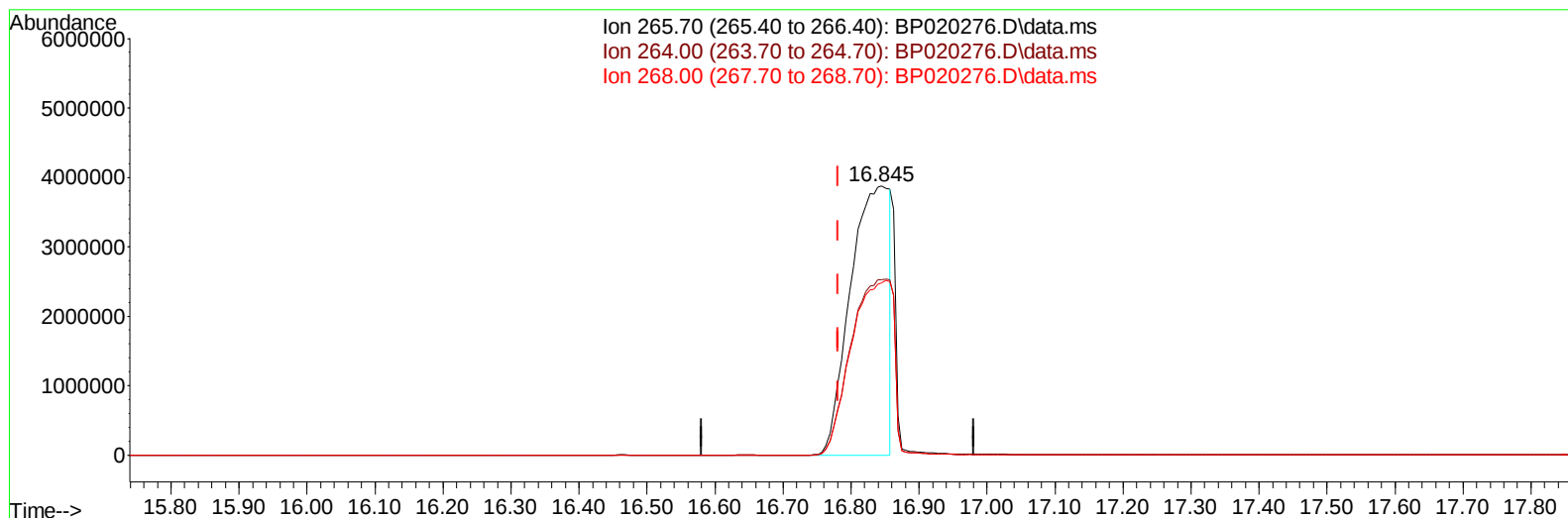
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TIC: BP020276.D\data.ms

(71) Pentachlorophenol (C)

16.845min (+ 0.065) 4115.55 ng/u1

response 15464148

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	62.20	65.15
268.00	63.00	64.05
0.00	0.00	0.00

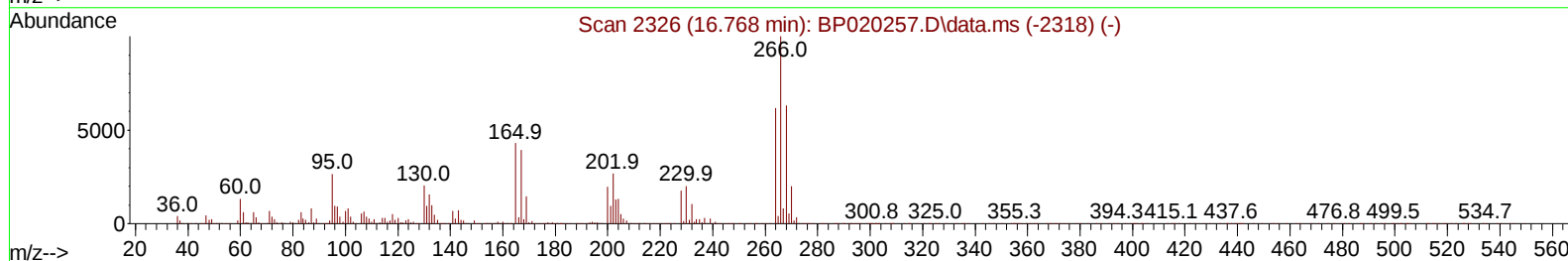
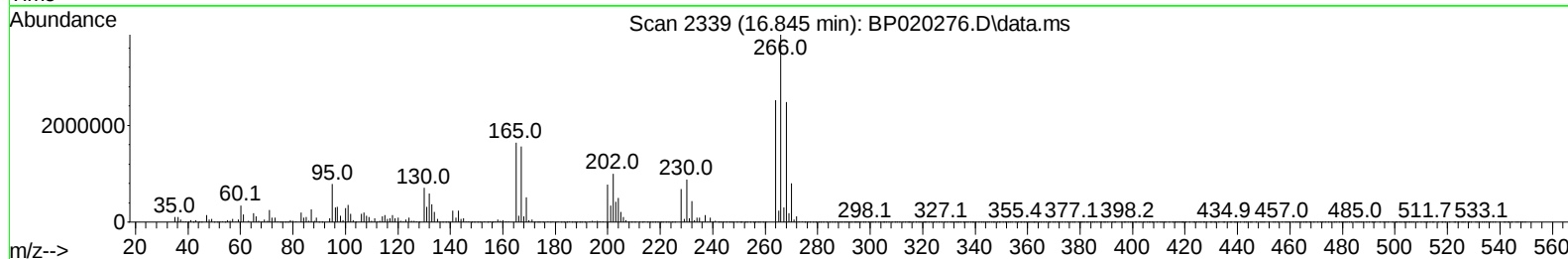
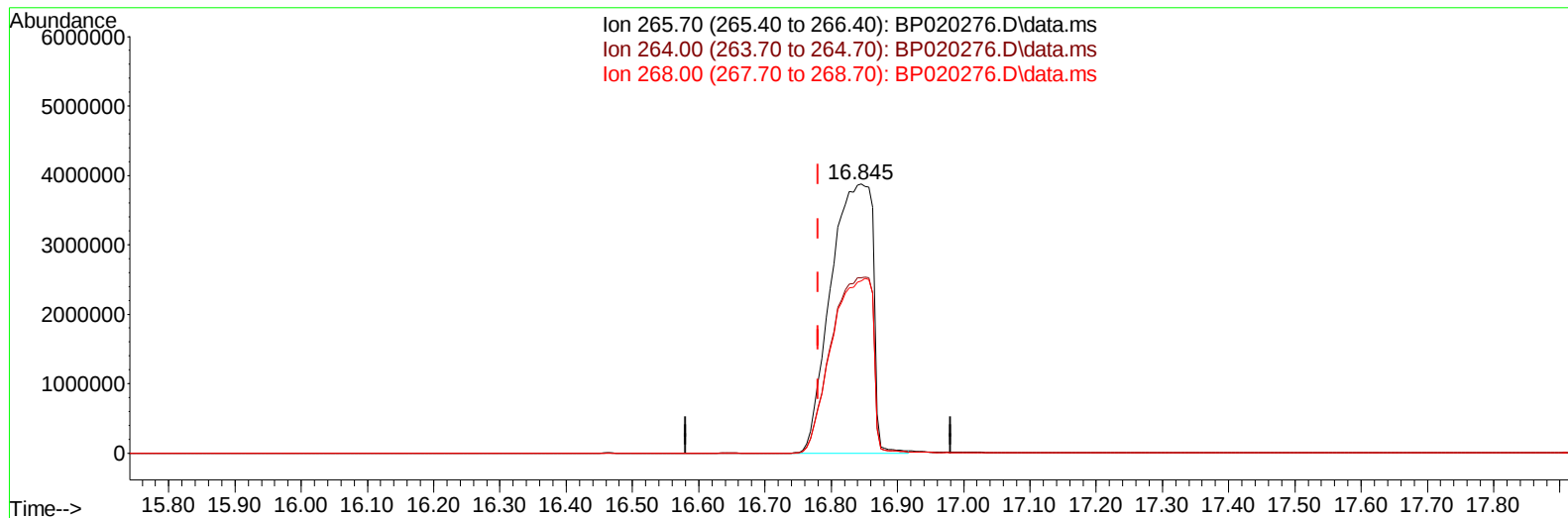
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 ALS Vial : 44 Sample Multiplier: 1

Instrument :
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TIC: BP020276.D\data.ms

(71) Pentachlorophenol (C)

16.845min (+ 0.065) 4539.60 ng/ul m

response 17057480

Ion	Exp%	Act%
265.70	100.00	100.00
264.00	62.20	65.15
268.00	63.00	64.05
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	78456	20.000	ng/ul	0.00
20) Naphthalene-d8	10.404	136	316444	20.000	ng/ul	# 0.02
38) Acenaphthene-d10	14.298	164	250084	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.157	188	495512	20.000	ng/ul	0.02
79) Chrysene-d12	21.639	240	504523	20.000	ng/ul	0.00
88) Perylene-d12	25.010	264	583153	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.217	96	8932	4.808	ng/uL	0.00
4) Pyridine-d5	3.622	84	111428	20.643	ng/ul	0.00
7) Phenol-d5	6.816	99	199172	29.561	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.987	67	411529	100.546	ng/ul	0.02
11) 2-Chlorophenol-d4	7.163	132	125607	26.078	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	121473	22.300	ng/ul	0.02
21) Nitrobenzene-d5	8.793	128	78946	33.316	ng/ul	0.01
24) 2-Nitrophenol-d4	9.504	143	71221	26.246	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.046	165	127172	25.682	ng/ul	0.02
31) 4-Chloroaniline-d4	10.563	131	64207	9.264	ng/ul	0.01
46) Dimethylphthalate-d6	13.710	166	404651	22.157	ng/ul	0.00
49) Acenaphthylene-d8	13.981	160	469954	23.295	ng/ul	0.00
54) 4-Nitrophenol-d4	14.587	143	64734m	23.063	ng/ul	0.04
60) Fluorene-d10	15.322	176	360851	23.689	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.487	200	21737	6.913	ng/ul	0.00
73) Anthracene-d10	17.257	188	576010	26.871	ng/ul	0.01
81) Pyrene-d10	19.645	212	707076	23.975	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.786	264	706283	25.049	ng/ul	0.01
Target Compounds						
					Qvalue	
71) Pentachlorophenol	16.845	266	17057480m	4539.597	ng/ul	

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

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