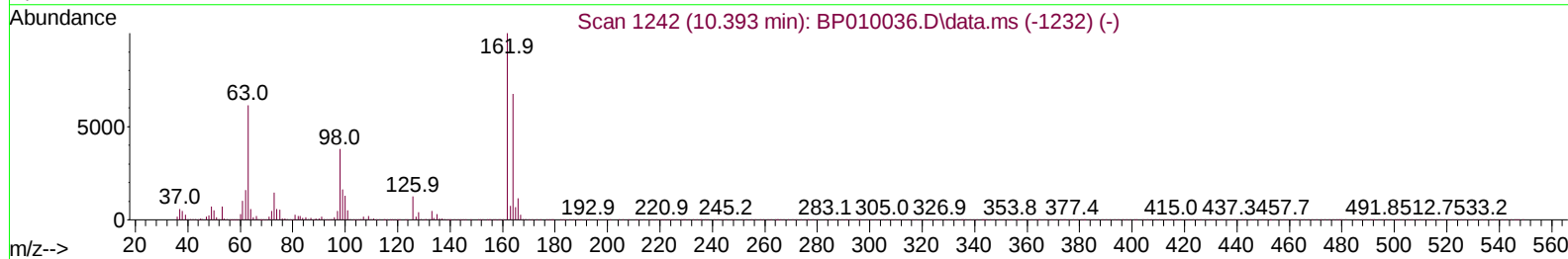
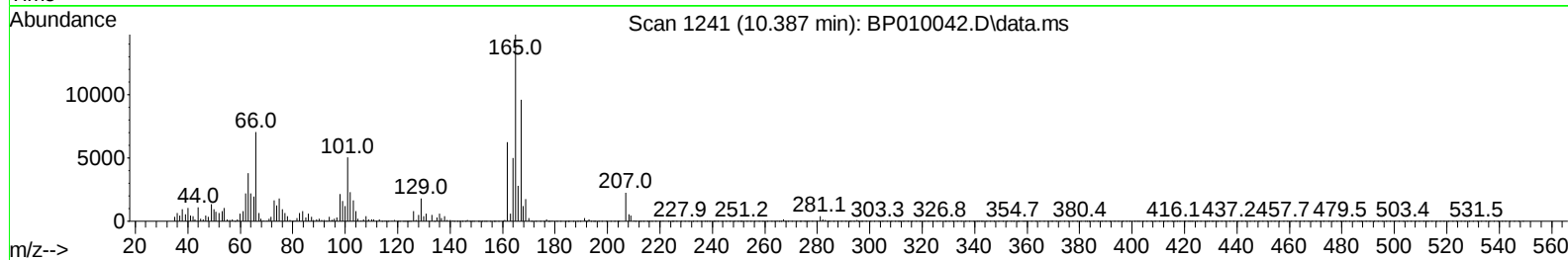
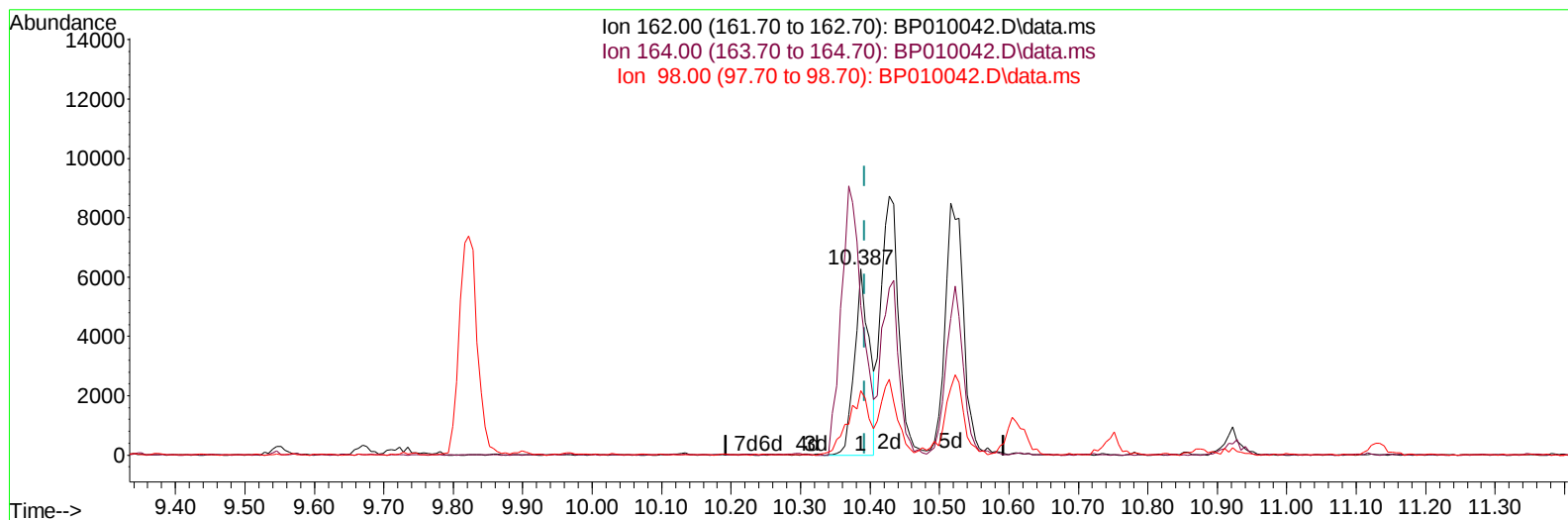


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
Data File : BP010042.D
Acq On : 22 Apr 2022 07:20
Operator : CG/JU
Sample : N2490-03
Misc :
ALS Vial : 38 Sample Multiplier: 1

Quant Time: Apr 22 11:20:16 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration



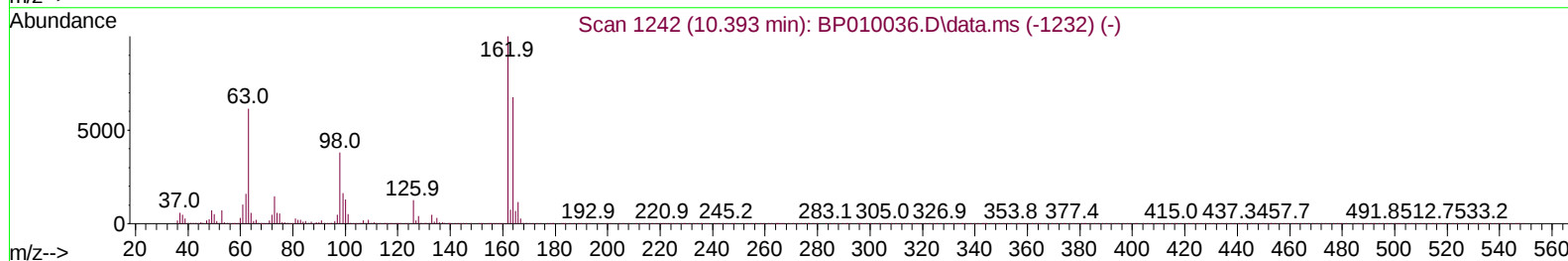
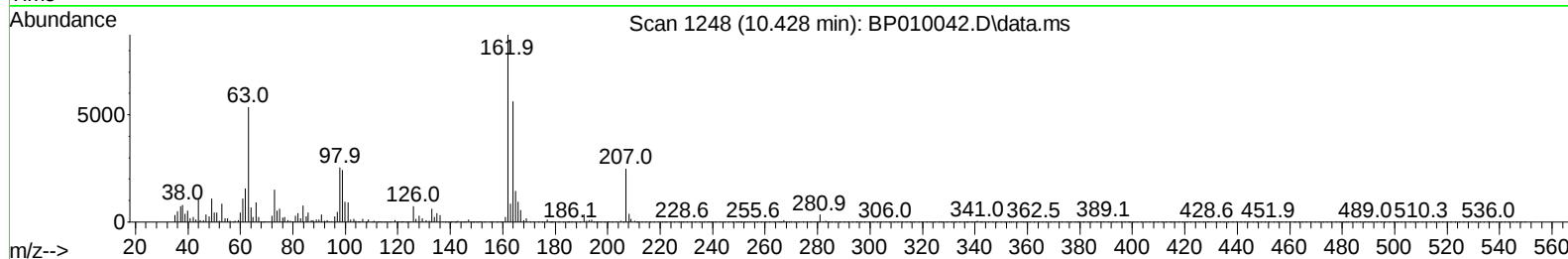
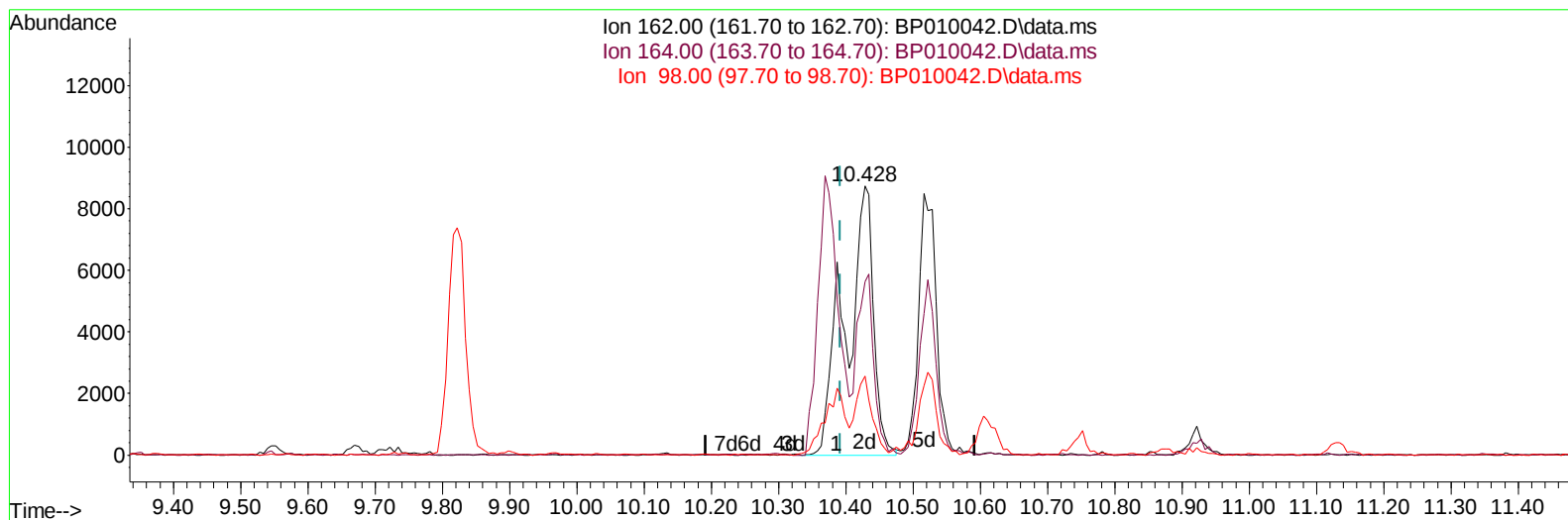
TIC: BP010042.D\data.ms

(29) 2,4-Dichlorophenol (C)**10.387min (-0.006) 0.88 ng/u1****response 9217**

Ion	Exp%	Act%
162.00	100.00	100.00
164.00	46.30	80.22#
98.00	38.80	34.68
0.00	0.00	0.00

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 Misc :
 ALS Vial : 38 Sample Multiplier: 1

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TIC: BP010042.D\data.ms

(29) 2,4-Dichlorophenol (C)**10.428min (+ 0.036) 2.37 ng/ul m****response 24806**

Ion	Exp%	Act%
162.00	100.00	100.00
164.00	46.30	64.55#
98.00	38.80	29.29#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042122\
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 Acq On : 22 Apr 2022 07:20
 Operator : CG/JU
 Sample : N2490-03
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Quant Time: Apr 22 11:20:16 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.940	152	135862	20.000	ng/ul	0.00
20) Naphthalene-d8	10.746	136	643997	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	473955	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.322	188	1046494	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.404	240	939884	20.000	ng/ul	# 0.00
88) Perylene-d12	23.857	264	887020	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.370	96	17605	5.651	ng/uL	0.00
4) Pyridine-d5	3.793	84	51161	5.837	ng/ul	0.00
7) Phenol-d5	7.099	99	75394	6.101	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.264	67	243584	32.849	ng/ul	0.00
11) 2-Chlorophenol-d4	7.469	132	222379	24.298	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	160480	15.674	ng/ul	-0.01
21) Nitrobenzene-d5	9.099	128	158322	30.910	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	167246	29.898	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	283126	25.999	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	347149	22.232	ng/ul	0.00
46) Dimethylphthalate-d6	13.975	166	1308373	35.488	ng/ul	0.00
49) Acenaphthylene-d8	14.263	160	1383261	31.000	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	36619m	5.395	ng/ul	0.00
60) Fluorene-d10	15.563	176	1129666	35.620	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	205919	31.119	ng/ul	0.00
73) Anthracene-d10	17.422	188	1950651	38.529	ng/ul	0.00
81) Pyrene-d10	19.651	212	2302775	43.082	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.698	264	1820754	39.154	ng/ul	0.00
Target Compounds						
					Qvalue	
12) 2-Chlorophenol	7.499	128	31038	3.362	ng/ul	95
22) Nitrobenzene	9.140	77	15447	1.231	ng/ul#	72
29) 2,4-Dichlorophenol	10.428	162	24806m	2.370	ng/ul	
39) 1,2,4,5-Tetrachloroben...	12.769	216	25008m	1.708	ng/ul	
75) 1,2,3,4-Tetrachloroben...	13.375	216	72619	4.753	ng/ul#	81

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

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