

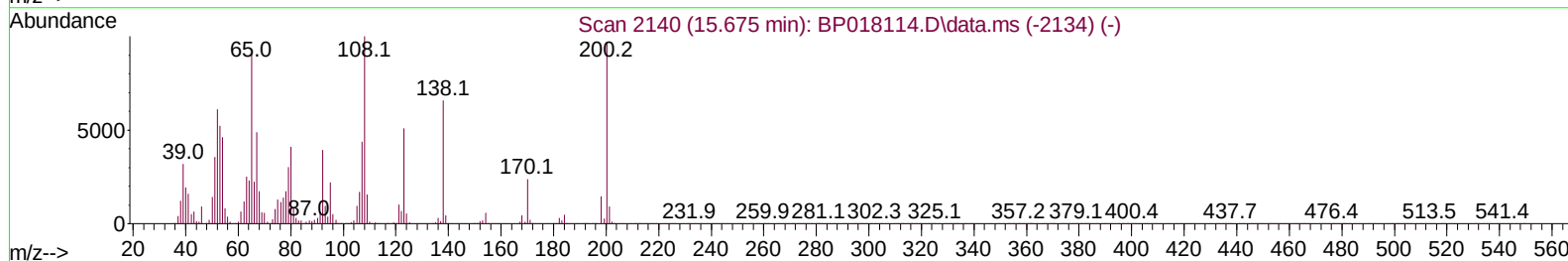
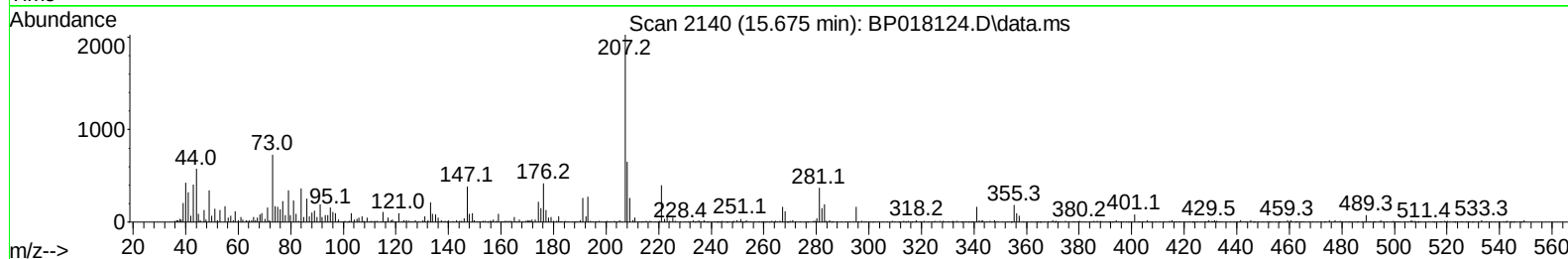
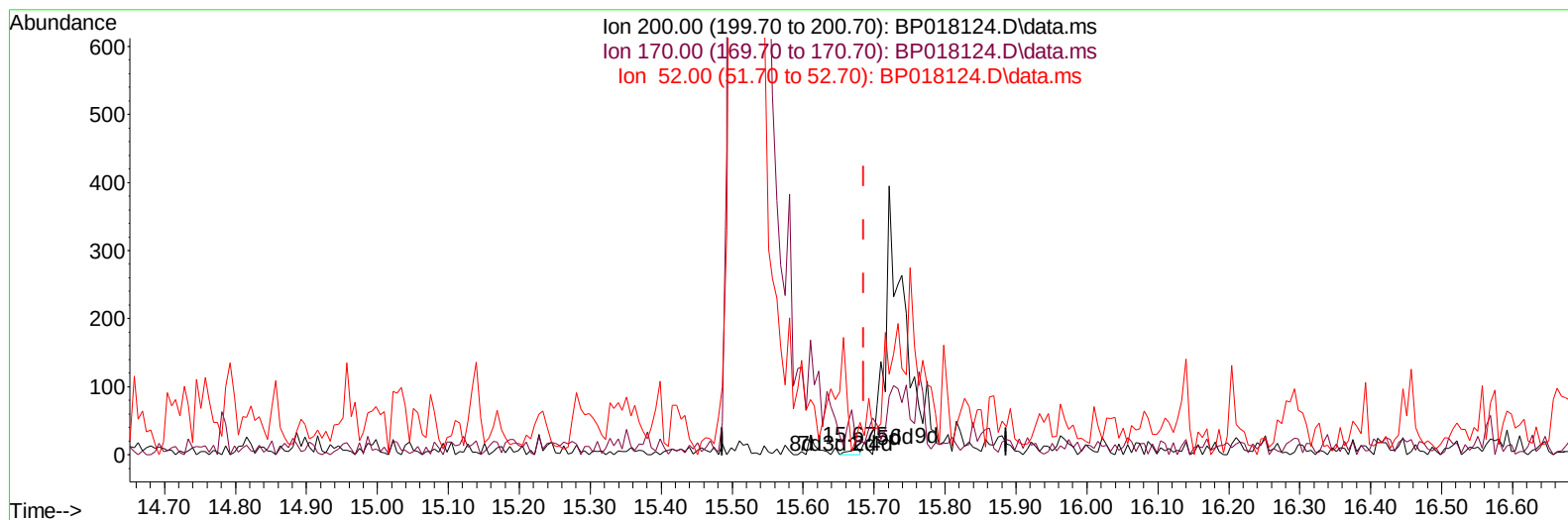
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018124.D
 Acq On : 13 Nov 2023 02:47
 Operator : MA/JU
 Sample : 05158-02
 Misc :
 ALS Vial : 105 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH363

Manual IntegrationsAPPROVED

Quant Time: Nov 13 03:51:04 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 10 21:55:18 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/14/2023
 Supervised By :mohammad ahmed 11/15/2023



TIC: BP018124.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.675min (-0.012) 0.00 ng/u1

response 11

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	269.23#
52.00	52.00	184.62#
0.00	0.00	0.00

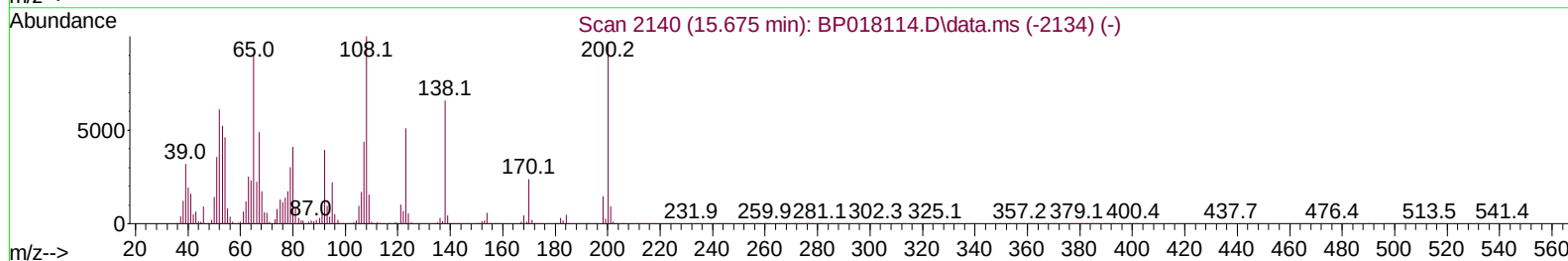
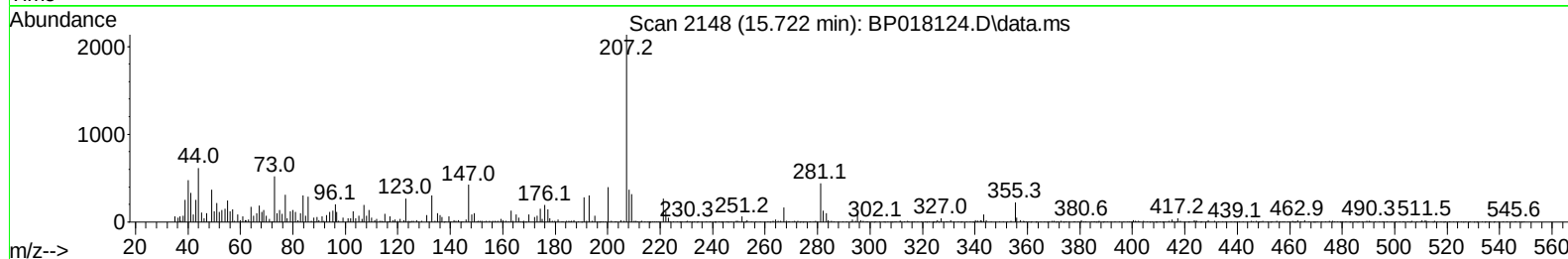
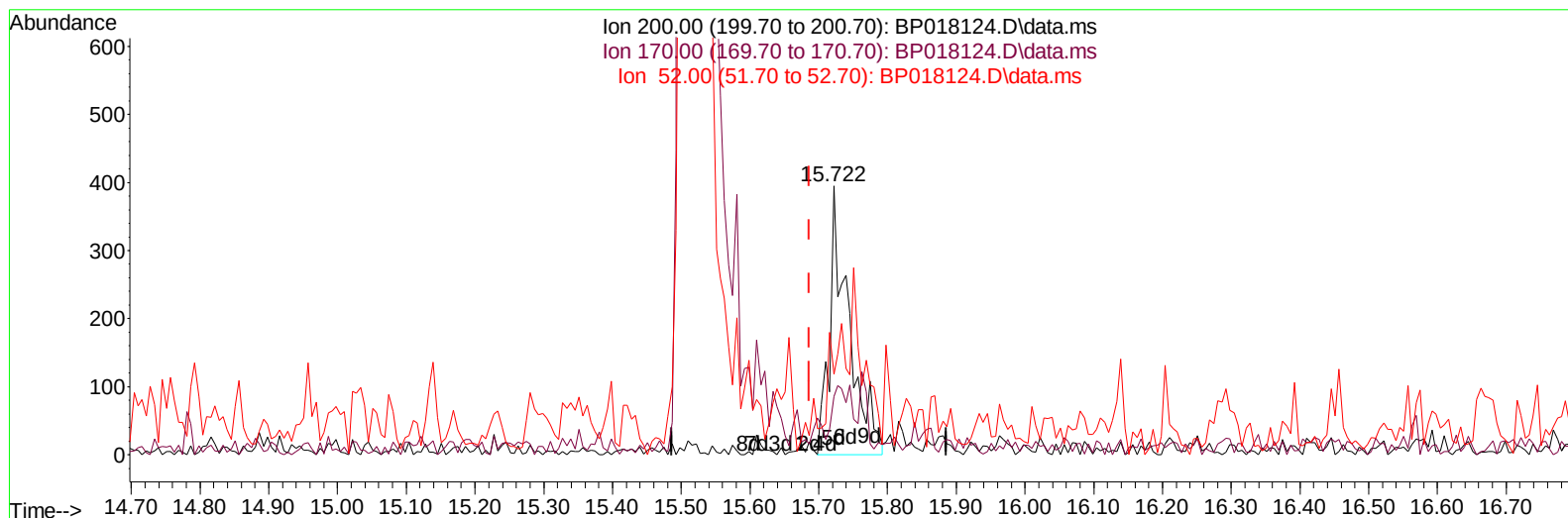
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.722min (+ 0.035) 0.21 ng/ul m

response 768

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	21.77
52.00	52.00	29.87#
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	101280	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.657	136	431905	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.516	164	263671	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.298	188	496233	20.000	ng/ul	-0.01
79) Chrysene-d12	21.427	240	402875	20.000	ng/ul	0.00
88) Perylene-d12	23.927	264	467890	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	6656	2.318	ng/uL	0.00
4) Pyridine-d5	3.669	84	43783	5.849	ng/ul	0.00
7) Phenol-d5	7.022	99	29739	3.041	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	97790	17.018	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	87523	11.303	ng/ul	0.00
15) 4-Methylphenol-d8	8.575	113	60221	7.516	ng/ul	0.00
21) Nitrobenzene-d5	9.051	128	56597	14.545	ng/ul	0.00
24) 2-Nitrophenol-d4	9.775	143	11610	2.634	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.298	165	56677	7.213	ng/ul	0.00
31) 4-Chloroaniline-d4	10.845	131	151804	14.089	ng/ul	0.00
46) Dimethylphthalate-d6	13.939	166	401405	16.499	ng/ul	-0.01
49) Acenaphthylene-d8	14.216	160	417143	15.743	ng/ul	-0.01
54) 4-Nitrophenol-d4	0.000	143	0d	0.000	ng/ul	
60) Fluorene-d10	15.516	176	327892	16.325	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.722	200	768m	0.214	ng/ul	0.04
73) Anthracene-d10	17.398	188	496222	18.489	ng/ul	-0.01
81) Pyrene-d10	19.651	212	555437	20.640	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.757	264	537154	19.744	ng/ul	-0.02

Target Compounds Qvalue

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

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