

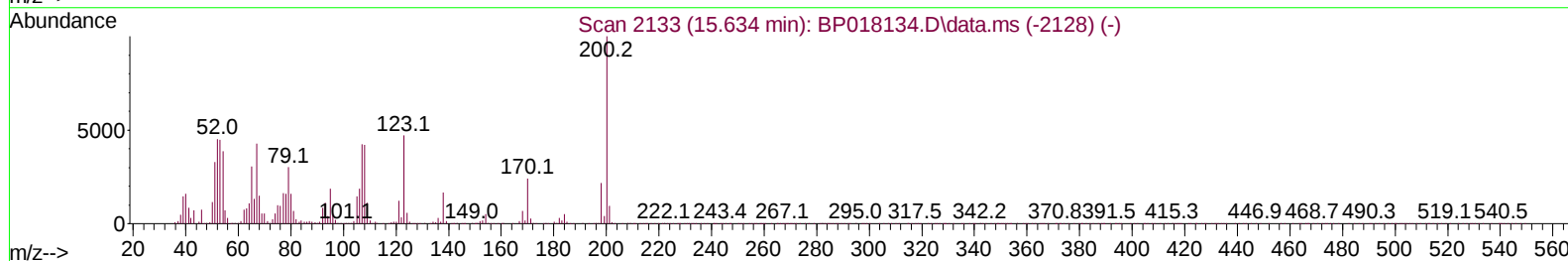
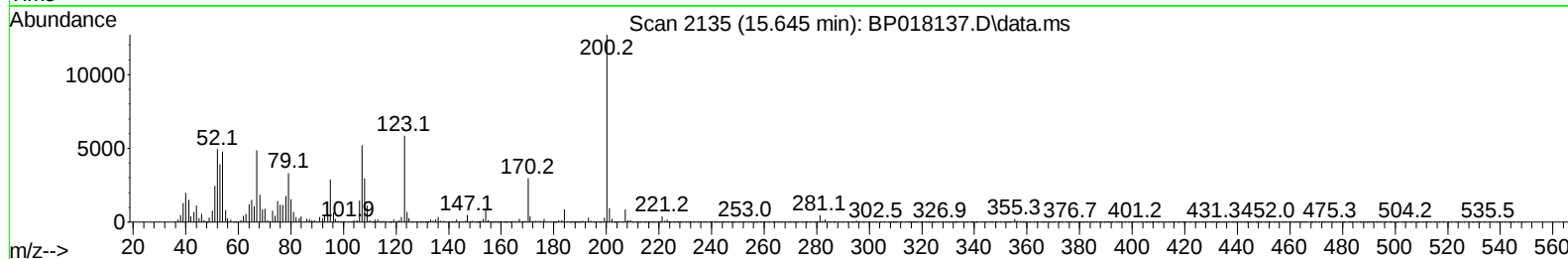
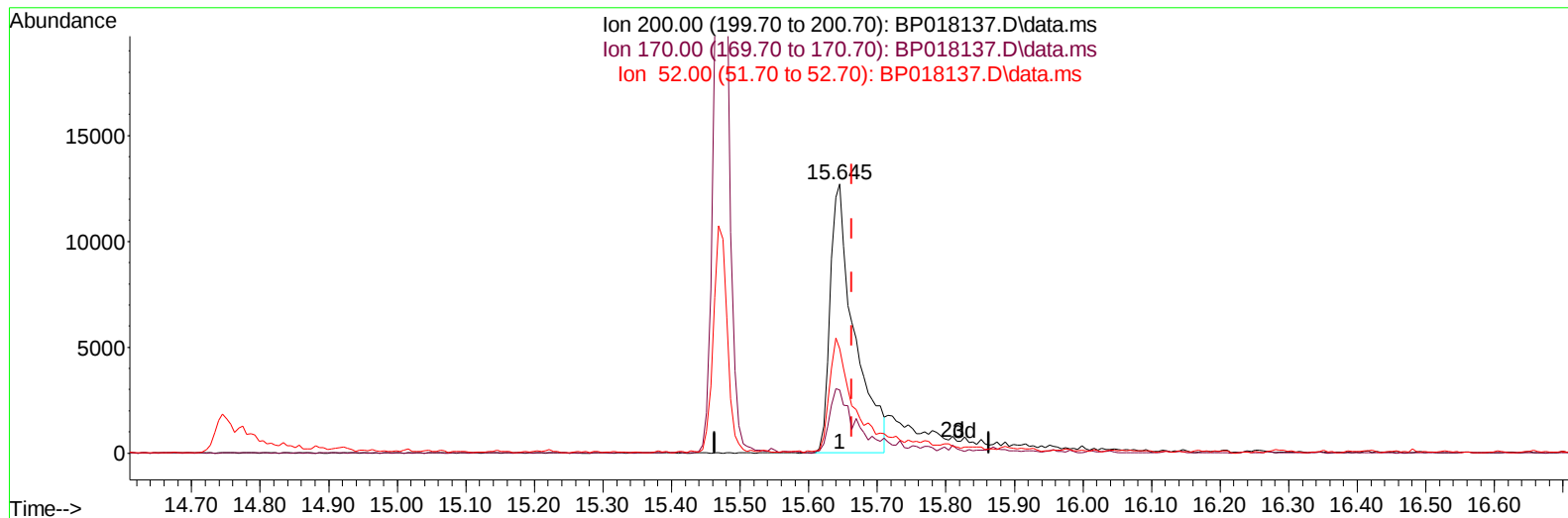
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111323\
 Data File : BP018137.D
 Acq On : 13 Nov 2023 18:33
 Operator : MA/JU
 Sample : PB156806BL
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK806

Manual IntegrationsAPPROVED

Quant Time: Nov 13 19:15:09 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 15:50:18 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/14/2023
 Supervised By :mohammad ahmed 11/16/2023



TIC: BP018137.D\data.ms

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

15.645min (-0.018) 8.84 ng/ul

response 30799

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	23.36
52.00	52.00	38.91#
0.00	0.00	0.00

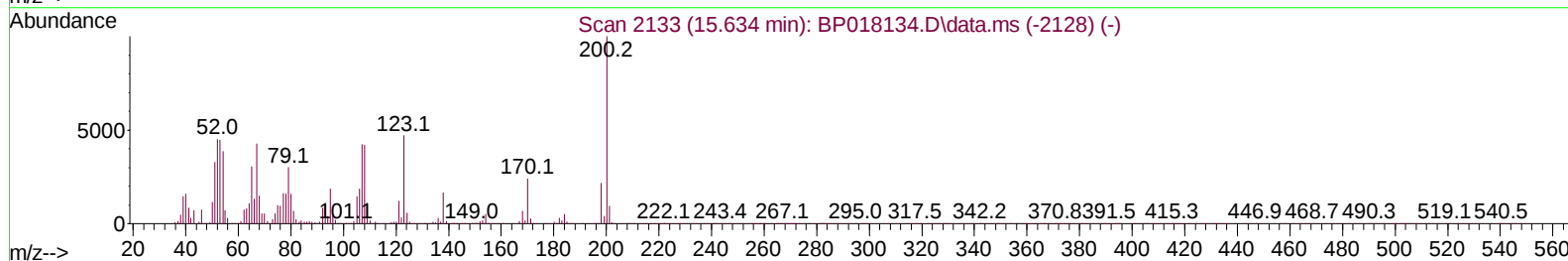
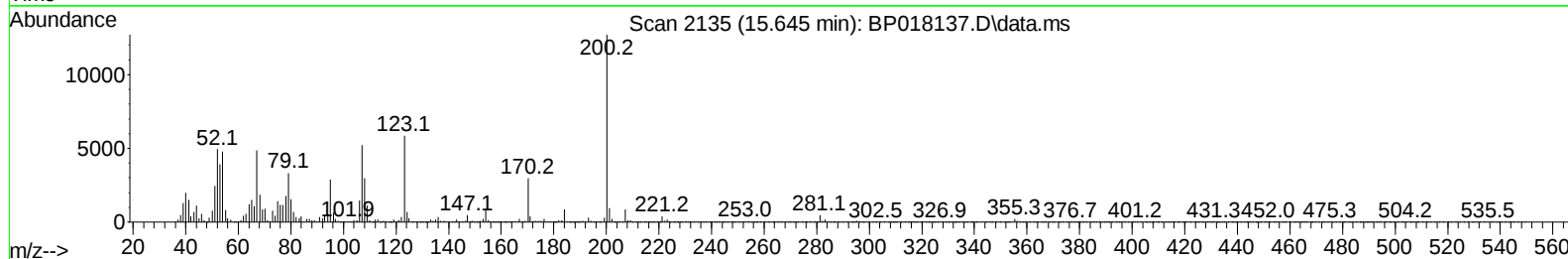
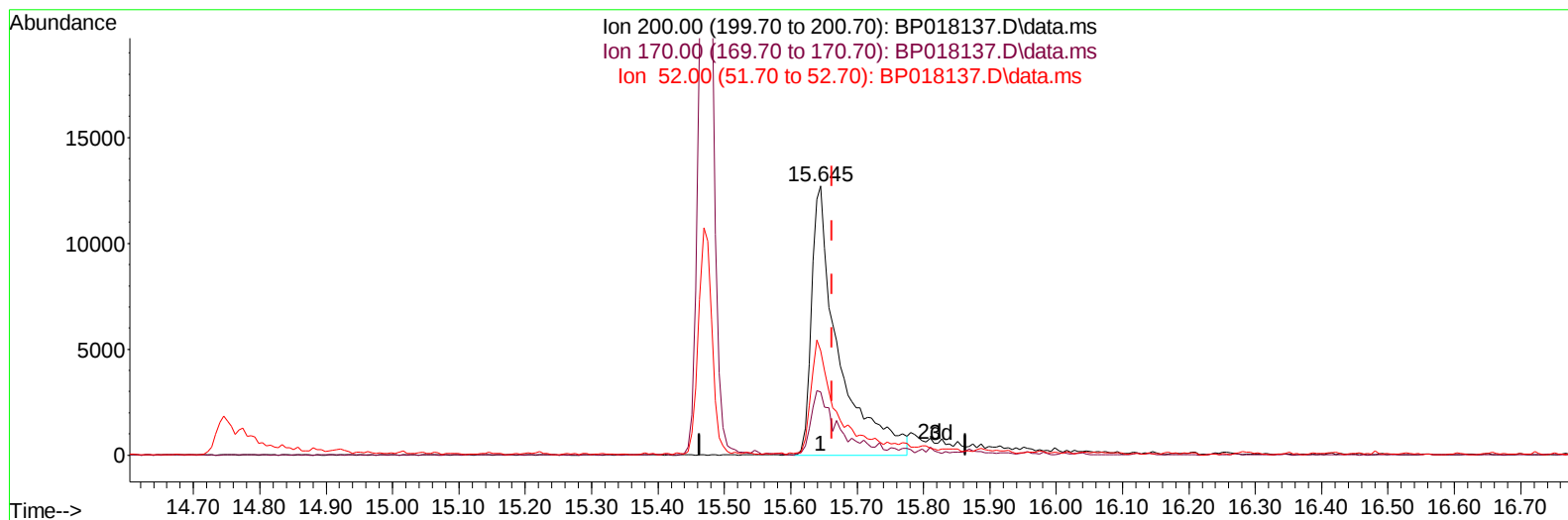
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TIC: BP018137.D\data.ms

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

15.645min (-0.018) 10.28 ng/ul m

response 35804

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	23.36
52.00	52.00	38.91#
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	7.793	152	84734	20.000	ng/ul	0.00
20) Naphthalene-d8	10.604	136	348479	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.469	164	183036	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.251	188	481931	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.380	240	423061	20.000	ng/ul	#-0.01
88) Perylene-d12	23.851	264	452872	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	13116	5.460	ng/uL	0.00
4) Pyridine-d5	3.605	84	95123	15.188	ng/ul	0.00
7) Phenol-d5	6.957	99	200102	24.461	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	134097	27.894	ng/ul	0.00
11) 2-Chlorophenol-d4	7.310	132	164104	25.332	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	141879	21.165	ng/ul	0.00
21) Nitrobenzene-d5	8.999	128	68387	21.782	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	90638	25.482	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	139075	21.938	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	162542	18.697	ng/ul	0.00
46) Dimethylphthalate-d6	13.892	166	520079	30.795	ng/ul	0.00
49) Acenaphthylene-d8	14.169	160	471990	25.659	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	20356	8.066	ng/ul	0.00
60) Fluorene-d10	15.475	176	380533	27.292	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	35804m	10.277	ng/ul	-0.02
73) Anthracene-d10	17.351	188	718985	27.585	ng/ul	0.00
81) Pyrene-d10	19.610	212	822823	29.117	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.686	264	739140	28.069	ng/ul	-0.02

Target Compounds Qvalue

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

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