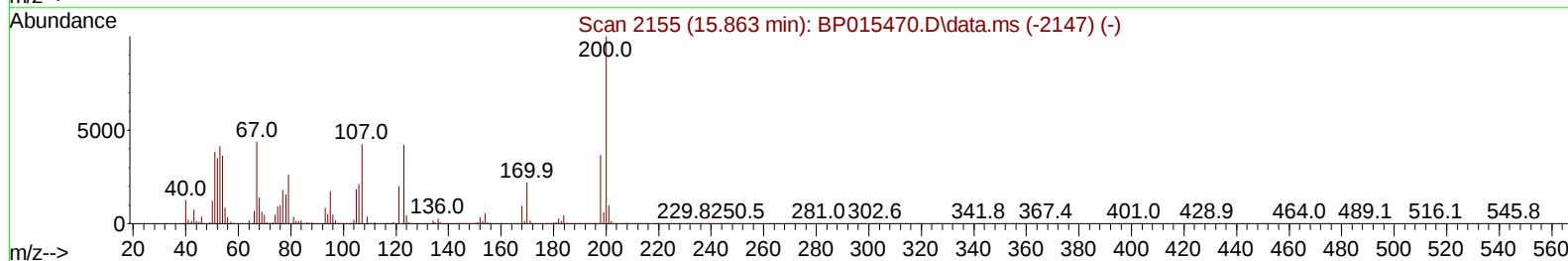
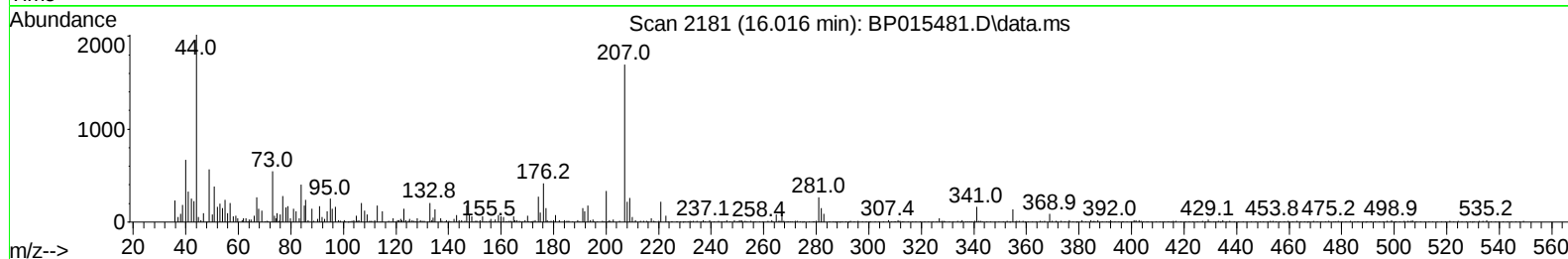
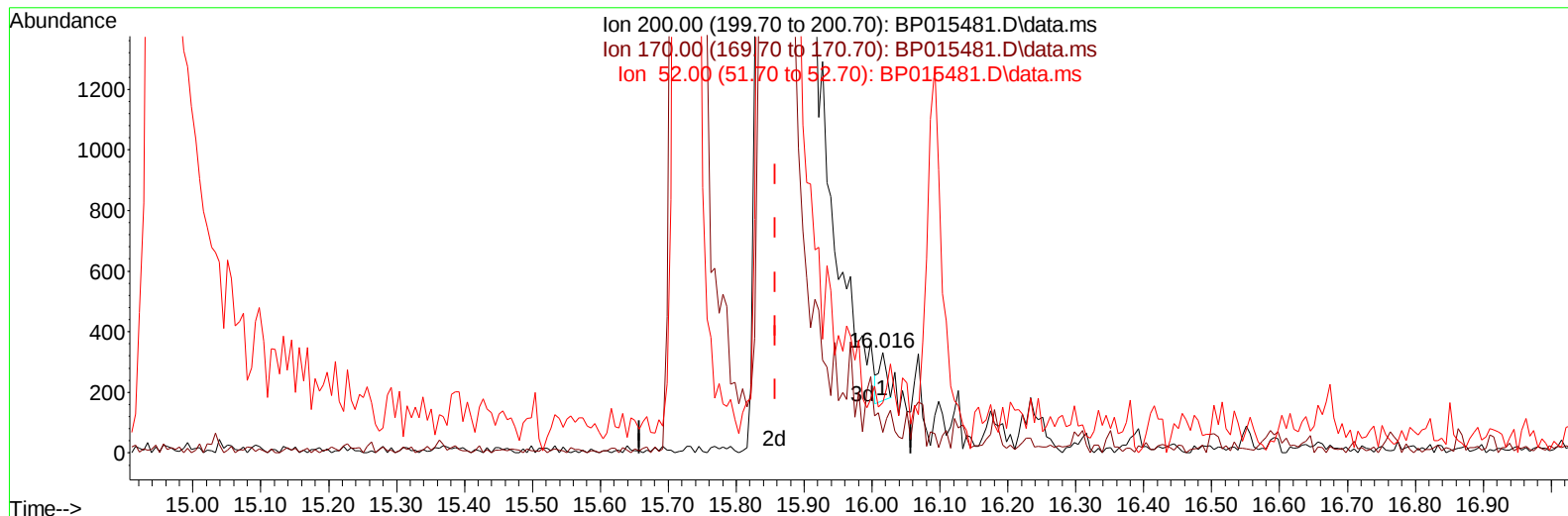


Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061223\
Data File : BP015481.D
Acq On : 12 Jun 2023 20:18
Operator : MA/JU
Sample : 03062-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jun 12 23:23:37 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 09 23:33:38 2023
Response via : Initial Calibration



TIC: BP015481.D\data.ms

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

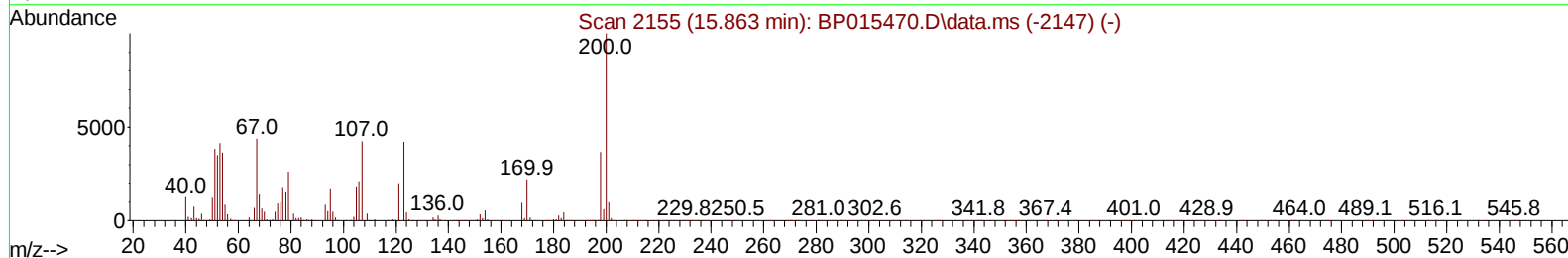
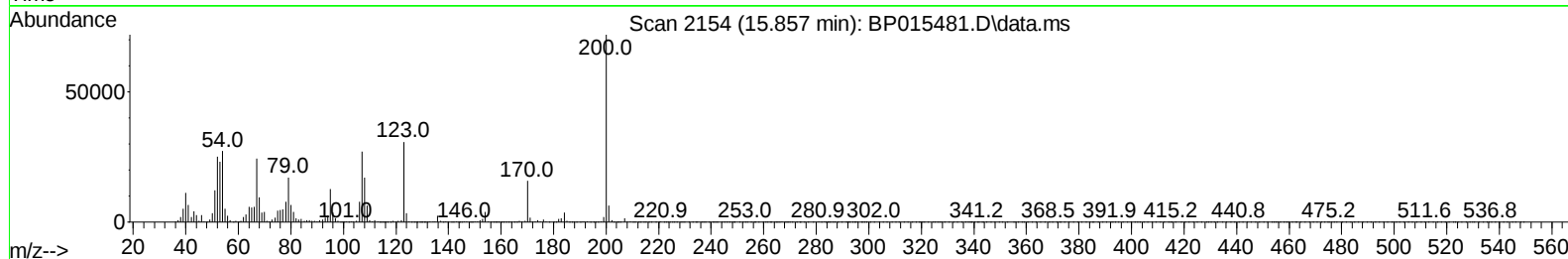
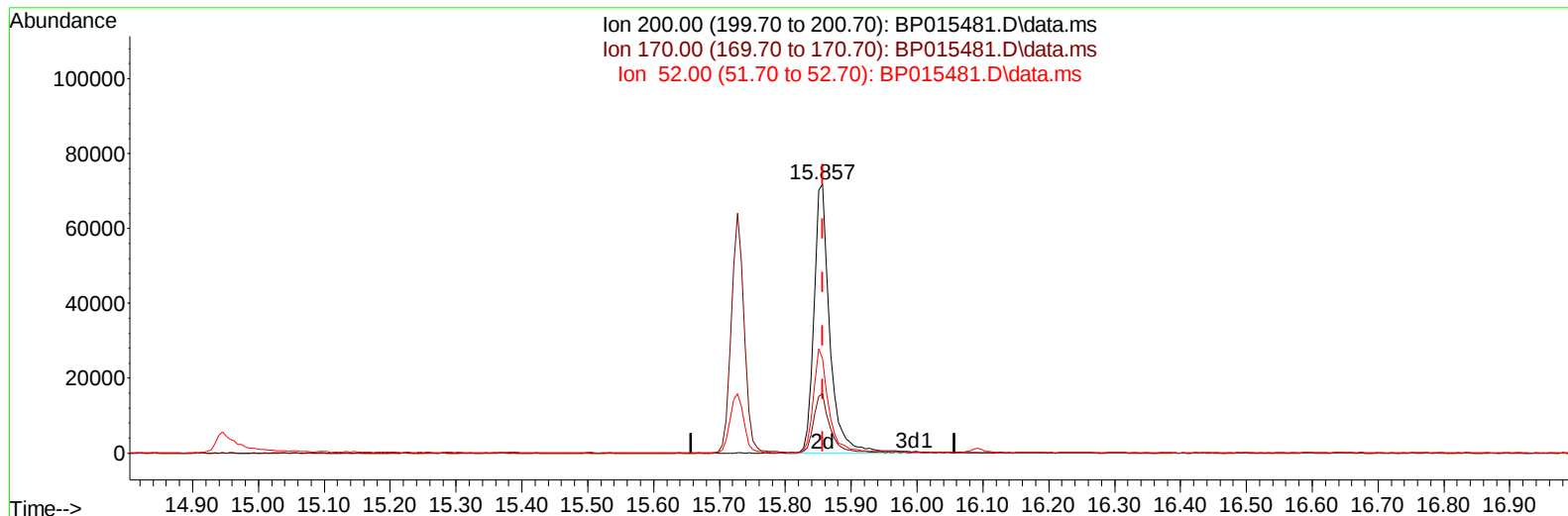
16.016min (+ 0.158) 0.02 ng/u1

response 123

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.30	22.71
52.00	46.20	48.97
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061223\
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Sample : 03062-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jun 12 23:23:37 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 09 23:33:38 2023
Response via : Initial Calibration



TIC: BP015481.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)**15.857min (-0.000) 19.55 ng/ul m****response 120291**

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.30	21.97
52.00	46.20	34.94#
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP061223\
 Data File : BP015481.D
 Acq On : 12 Jun 2023 20:18
 Operator : MA/JU
 Sample : 03062-10
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Quant Time: Jun 12 23:44:41 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Jun 09 23:33:38 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.092	152	120578	20.000	ng/ul	0.00
20) Naphthalene-d8	10.922	136	584737	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.733	164	417867	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.492	188	971323	20.000	ng/ul	0.00
79) Chrysene-d12	21.574	240	959214	20.000	ng/ul	0.00
88) Perylene-d12	24.115	264	1012024	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.410	96	15873	4.875	ng/uL	0.00
4) Pyridine-d5	3.852	84	175940	20.801	ng/ul	0.00
7) Phenol-d5	7.245	99	251544	22.644	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.416	67	164425	24.785	ng/ul	0.00
11) 2-Chlorophenol-d4	7.616	132	198956	24.703	ng/ul	0.00
15) 4-Methylphenol-d8	8.804	113	223994	24.569	ng/ul	0.00
21) Nitrobenzene-d5	9.269	128	106011	23.092	ng/ul	0.00
24) 2-Nitrophenol-d4	9.998	143	126683	24.164	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.539	165	215726	22.557	ng/ul	0.00
31) 4-Chloroaniline-d4	11.063	131	320333	23.336	ng/ul	0.00
46) Dimethylphthalate-d6	14.139	166	865452	26.265	ng/ul	0.00
49) Acenaphthylene-d8	14.427	160	935606	25.967	ng/ul	0.00
54) 4-Nitrophenol-d4	14.945	143	89518	15.170	ng/ul	0.00
60) Fluorene-d10	15.727	176	731732	26.334	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	120291m	19.549	ng/ul	0.00
73) Anthracene-d10	17.592	188	1178135	26.658	ng/ul	0.00
81) Pyrene-d10	19.809	212	1446612	28.178	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.950	264	1350480	26.595	ng/ul	0.00

Target Compounds	Qvalue

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

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