

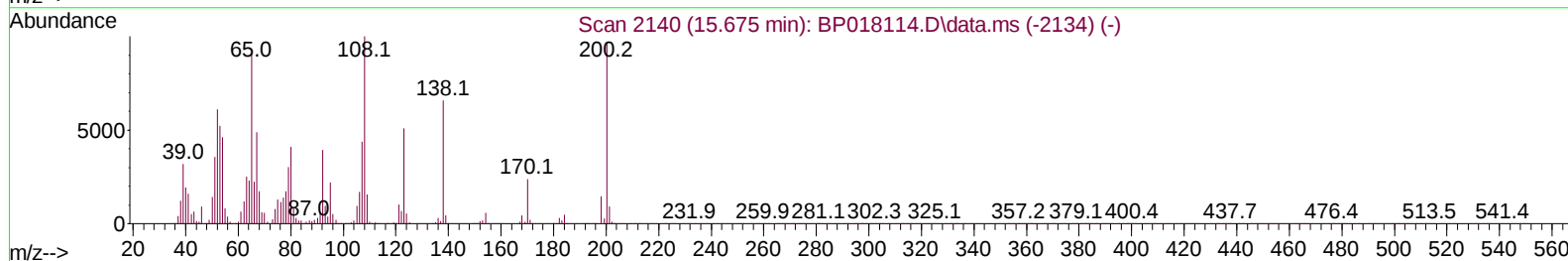
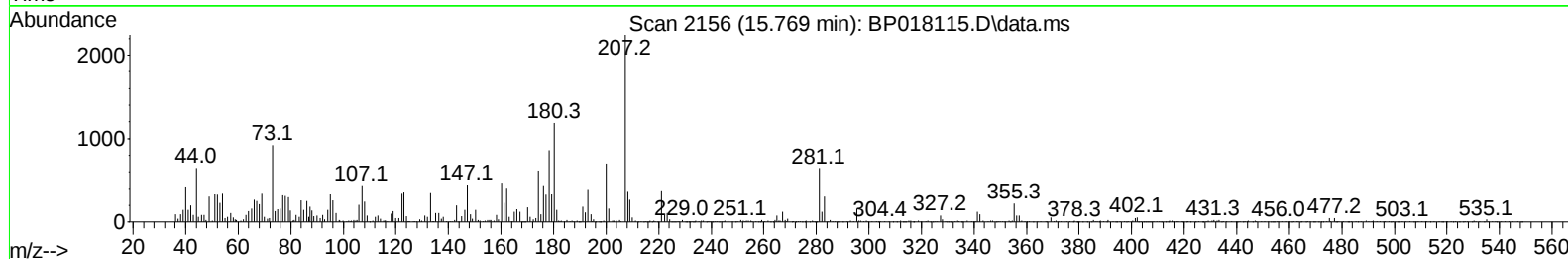
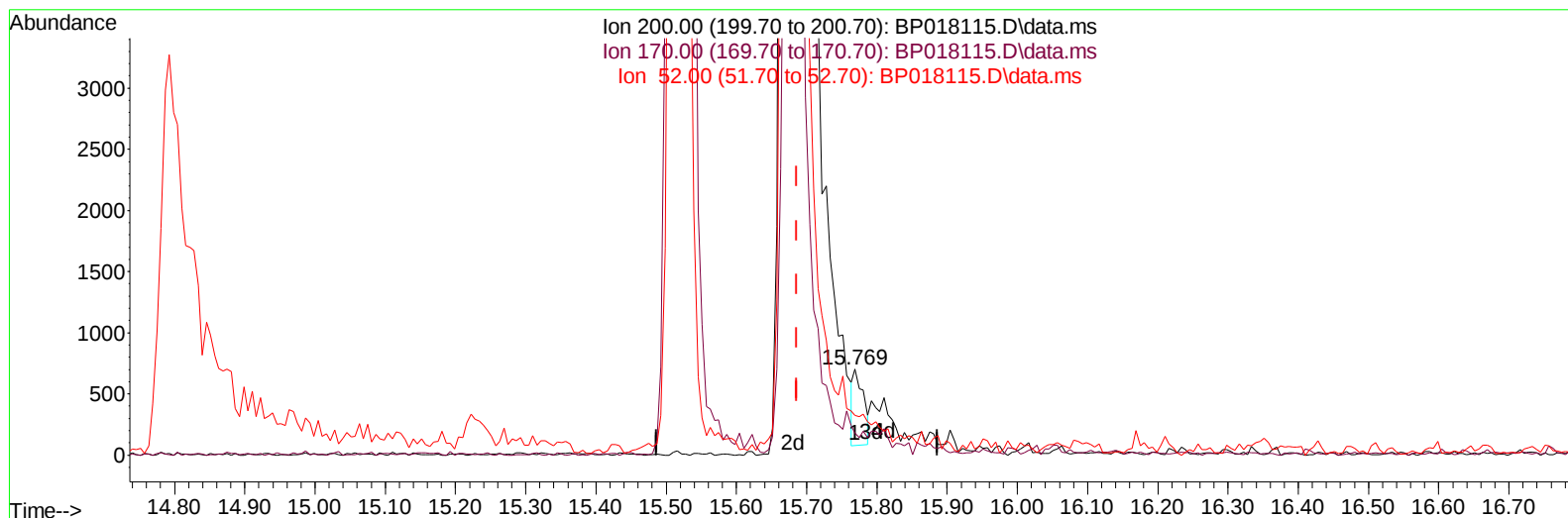
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111023\
 Data File : BP018115.D
 Acq On : 12 Nov 2023 21:42
 Operator : MA/JU
 Sample : PB156723BL
 Misc :
 ALS Vial : 96 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK723

Manual IntegrationsAPPROVED

Quant Time: Nov 13 03:05:46 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: BP018115.D\data.ms

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

15.769min (+ 0.083) 0.21 ng/ul

response 629

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	25.36
52.00	52.00	46.15
0.00	0.00	0.00

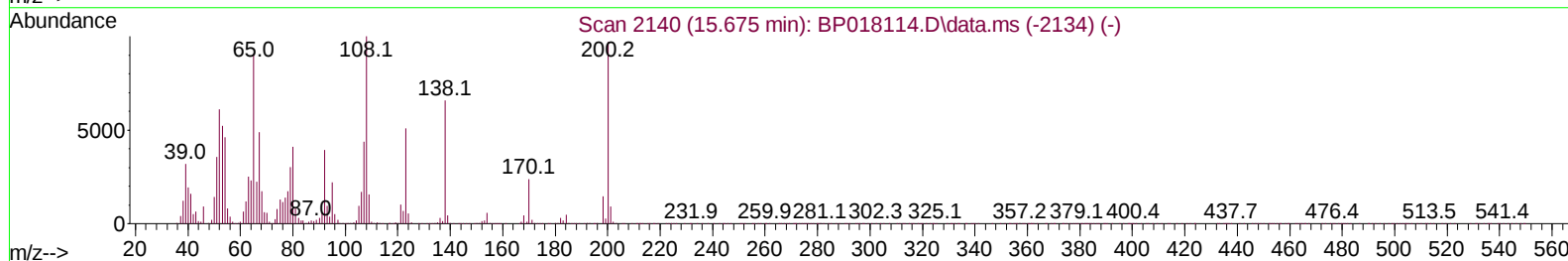
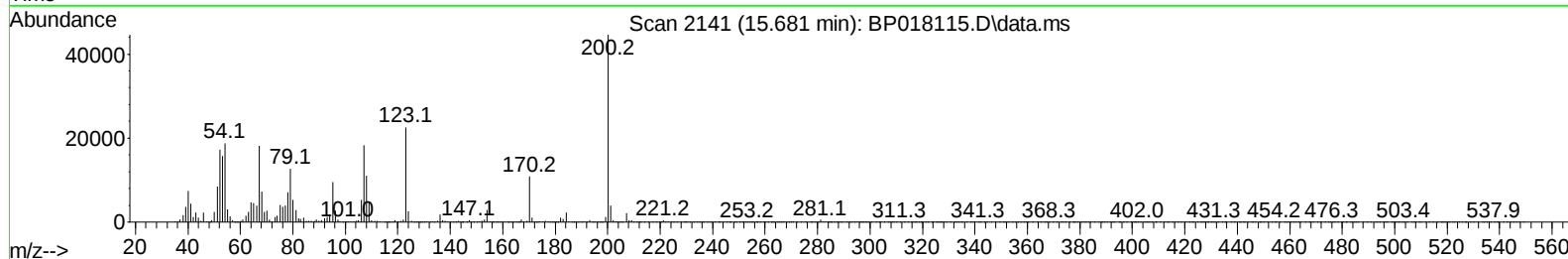
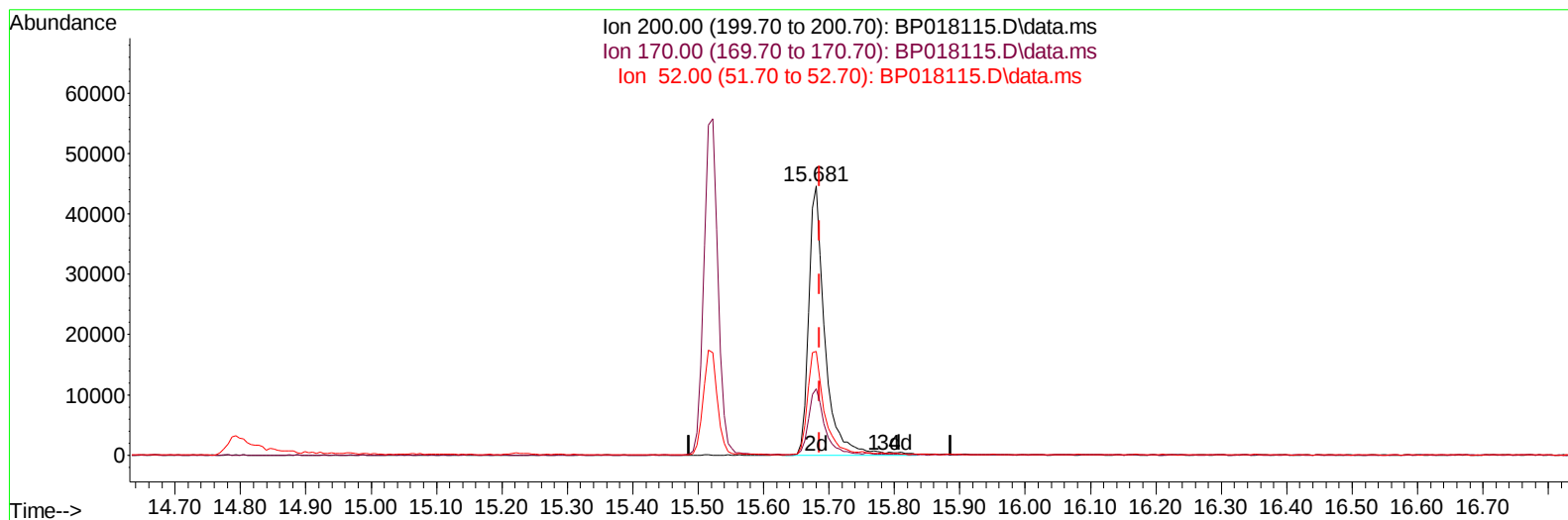
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TIC: BP018115.D\data.ms

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

15.681min (-0.006) 25.74 ng/ul m

response 76513

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	24.56
52.00	52.00	38.63#
0.00	0.00	0.00

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Manual IntegrationsAPPROVED

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Quant Time: Nov 13 03:09:00 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.840	152	75519	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.663	136	325201	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.516	164	208206	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.298	188	411194	20.000	ng/ul	-0.01
79) Chrysene-d12	21.427	240	280321	20.000	ng/ul	0.00
88) Perylene-d12	23.927	264	302210	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	17056	7.967	ng/uL	0.00
4) Pyridine-d5	3.658	84	219488	39.322	ng/ul	0.00
7) Phenol-d5	7.011	99	302369	41.472	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.193	67	188786	44.061	ng/ul	0.00
11) 2-Chlorophenol-d4	7.364	132	238064	41.233	ng/ul	-0.01
15) 4-Methylphenol-d8	8.563	113	250327	41.900	ng/ul	-0.01
21) Nitrobenzene-d5	9.046	128	121230	41.377	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.763	143	141500	42.630	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	237122	40.082	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.840	131	304880	37.579	ng/ul	-0.01
46) Dimethylphthalate-d6	13.940	166	807106	42.013	ng/ul	-0.01
49) Acenaphthylene-d8	14.216	160	857101	40.963	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.793	143	57820	20.142	ng/ul	0.01
60) Fluorene-d10	15.522	176	651289	41.063	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.681	200	76513m	25.739	ng/ul	0.00
73) Anthracene-d10	17.398	188	939552	42.248	ng/ul	-0.01
81) Pyrene-d10	19.651	212	955660	51.038	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.763	264	780566	44.419	ng/ul	-0.01

Target Compounds Qvalue

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

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