

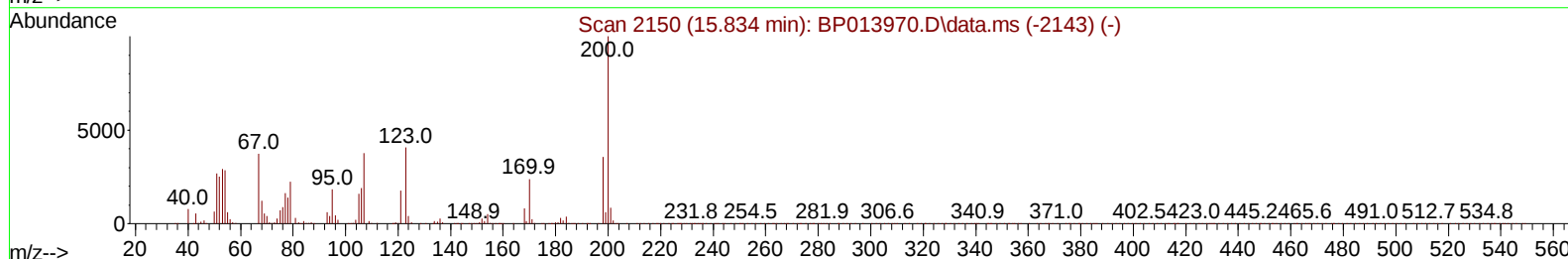
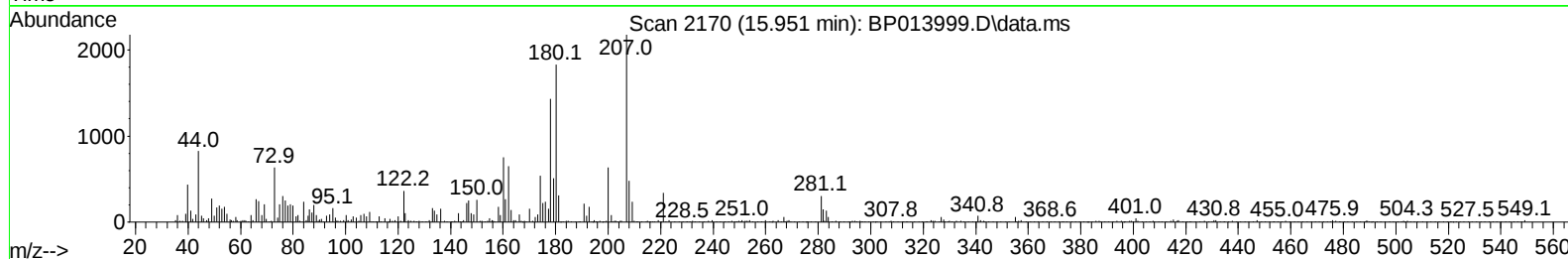
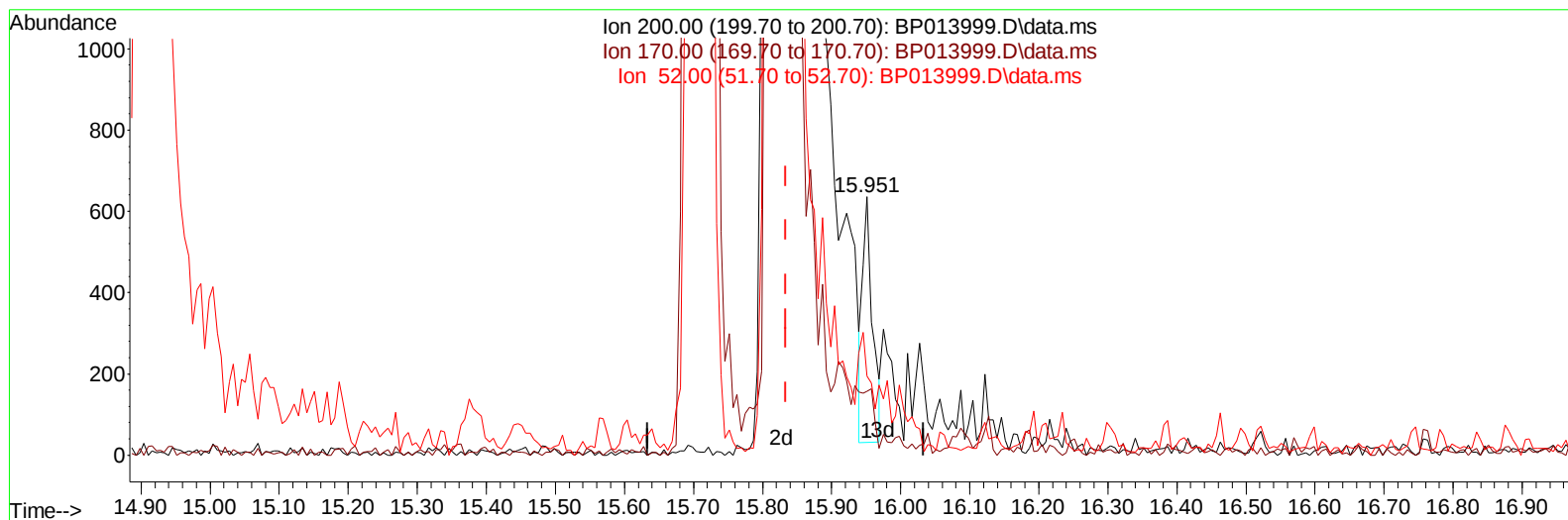
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
 Data File : BP013999.D
 Acq On : 09 Mar 2023 03:25
 Operator : CG/JU
 Sample : PB151195BL
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK195

Manual IntegrationsAPPROVED

Quant Time: Mar 09 04:33:33 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP030723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 09 01:10:12 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 03/09/2023
 Supervised By :Jagrut Upadhyay 03/09/2023



TIC: BP013999.D\data.ms

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

15.951min (+ 0.118) 0.08 ng/ul

response 608

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.80	24.69
52.00	37.50	30.50
0.00	0.00	0.00

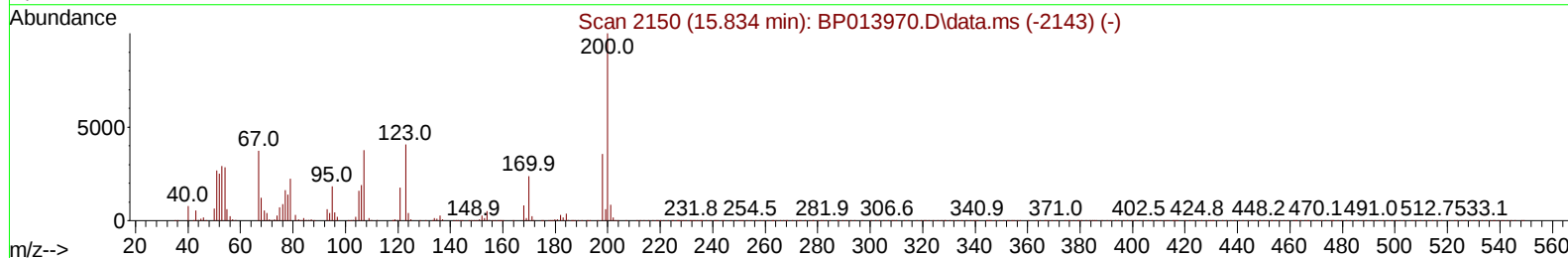
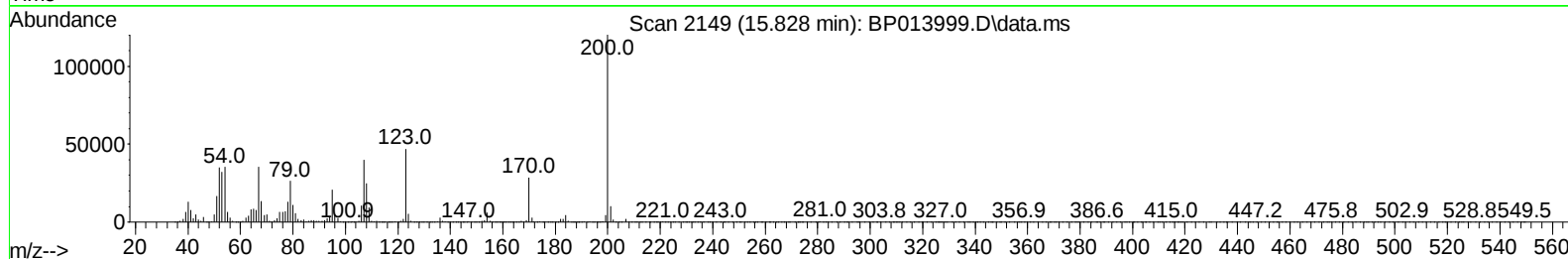
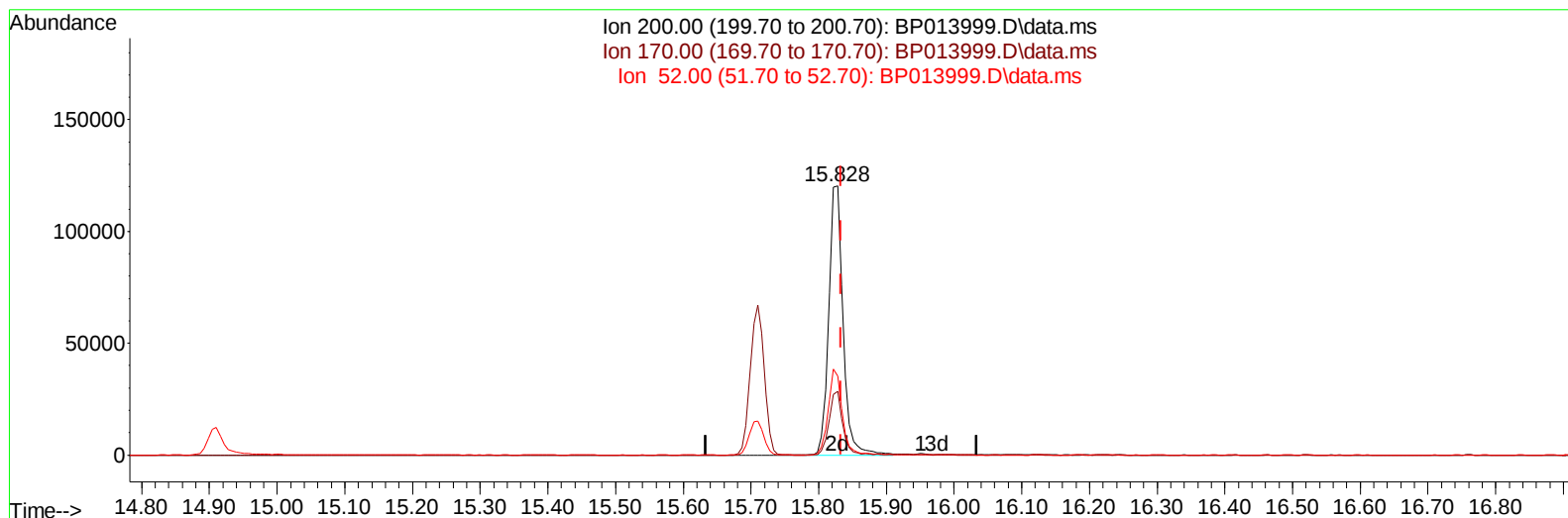
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TIC: BP013999.D\data.ms

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

15.828min (-0.006) 24.21 ng/ul m

response 176163

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.80	23.83
52.00	37.50	29.26#
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	8.081	152	134201	20.000	ng/ul	0.00
20) Naphthalene-d8	10.905	136	539340	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.716	164	358514	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.469	188	962650	20.000	ng/ul	0.00
79) Chrysene-d12	21.545	240	1171487	20.000	ng/ul	-0.01
88) Perylene-d12	24.086	264	1306285	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.440	96	19489	5.487	ng/uL	0.00
4) Pyridine-d5	3.870	84	266675	27.117	ng/ul	0.00
7) Phenol-d5	7.228	99	328835	27.767	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.399	67	195094	28.247	ng/ul	0.00
11) 2-Chlorophenol-d4	7.605	132	251616	27.886	ng/ul	0.00
15) 4-Methylphenol-d8	8.787	113	262299	27.115	ng/ul	0.00
21) Nitrobenzene-d5	9.252	128	125524	29.169	ng/ul	0.00
24) 2-Nitrophenol-d4	9.981	143	150019	29.122	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.522	165	263996	27.236	ng/ul	0.00
31) 4-Chloroaniline-d4	11.040	131	342924	26.734	ng/ul	0.00
46) Dimethylphthalate-d6	14.122	166	916009	30.106	ng/ul	0.00
49) Acenaphthylene-d8	14.410	160	910448	28.247	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	142834	26.639	ng/ul	0.00
60) Fluorene-d10	15.710	176	773650	30.002	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.828	200	176163m	24.212	ng/ul	0.00
73) Anthracene-d10	17.569	188	1366048	29.231	ng/ul	0.00
81) Pyrene-d10	19.792	212	1877285	29.055	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.921	264	2073984	29.848	ng/ul	0.00

Target Compounds Qvalue

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

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