

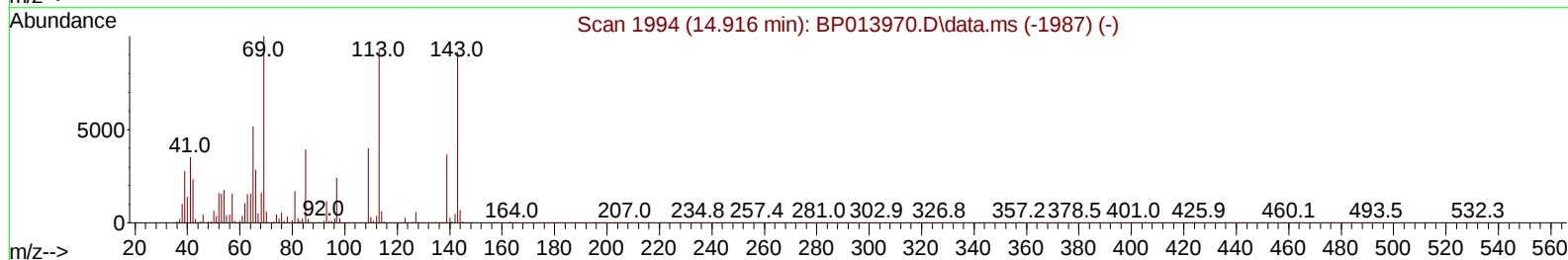
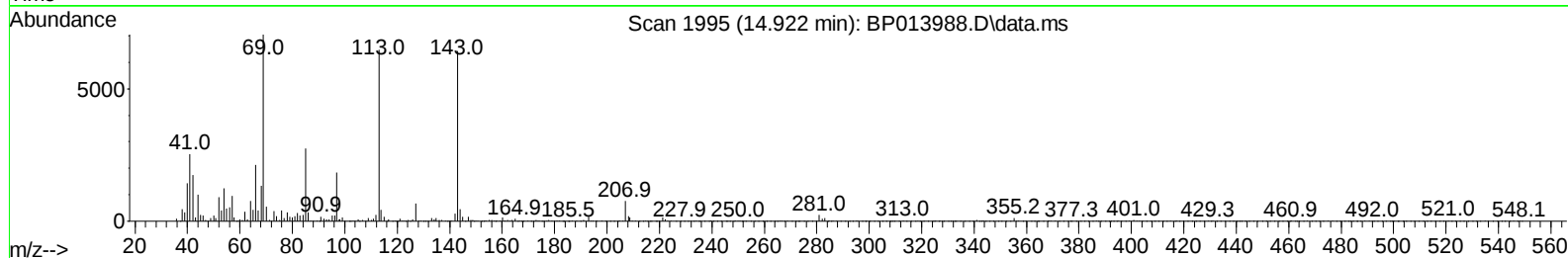
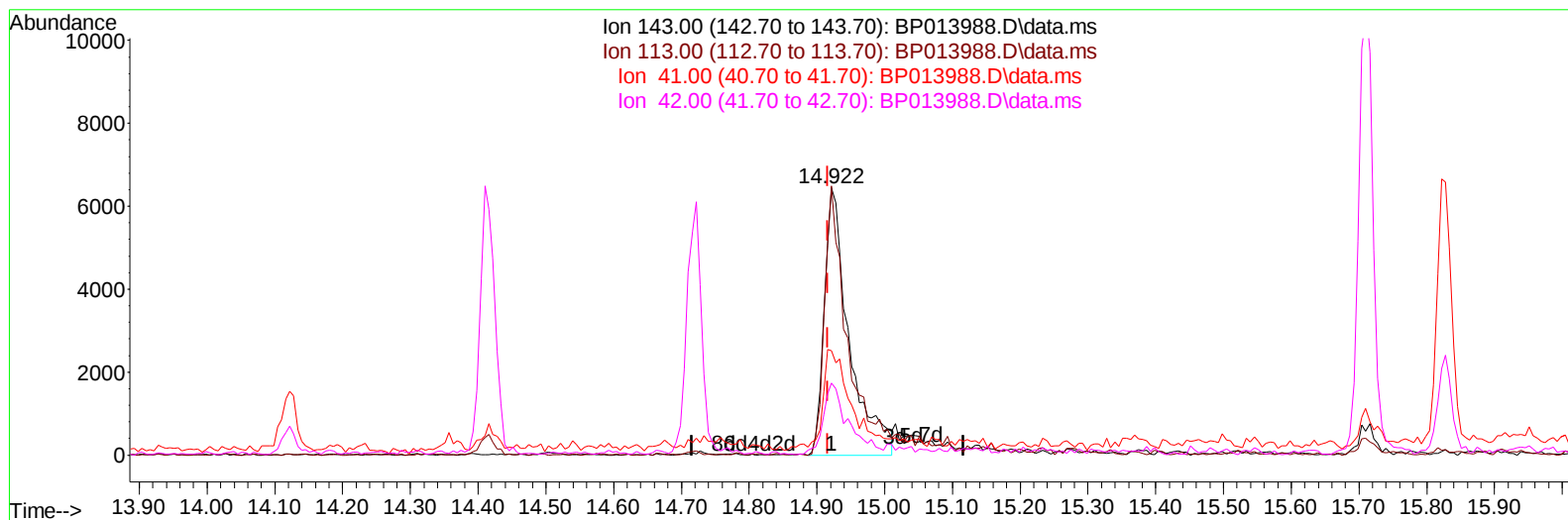
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP030823\
 Data File : BP013988.D
 Acq On : 08 Mar 2023 21:11
 Operator : CG/JU
 Sample : 01729-03
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0BY7

Manual IntegrationsAPPROVED

Quant Time: Mar 09 01:17:22 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: BP013988.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.922min (+ 0.006) 4.14 ng/ul m

response 16457

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	103.10	100.73
41.00	40.20	39.18
42.00	28.50	27.02

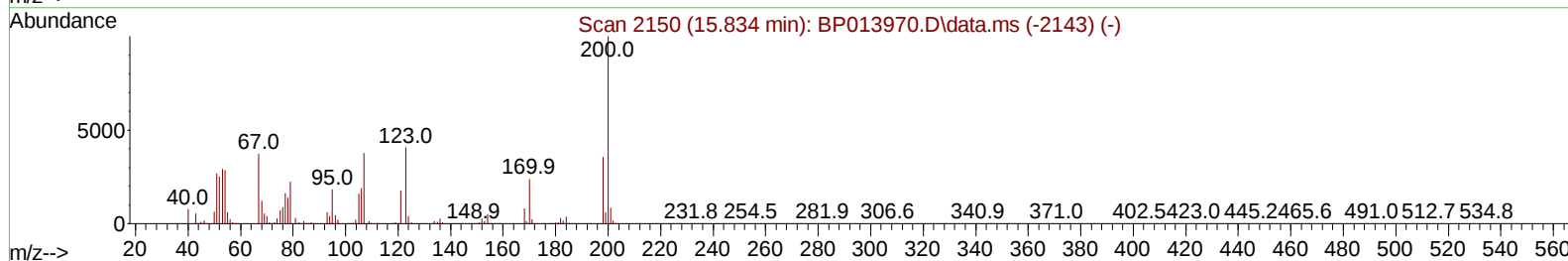
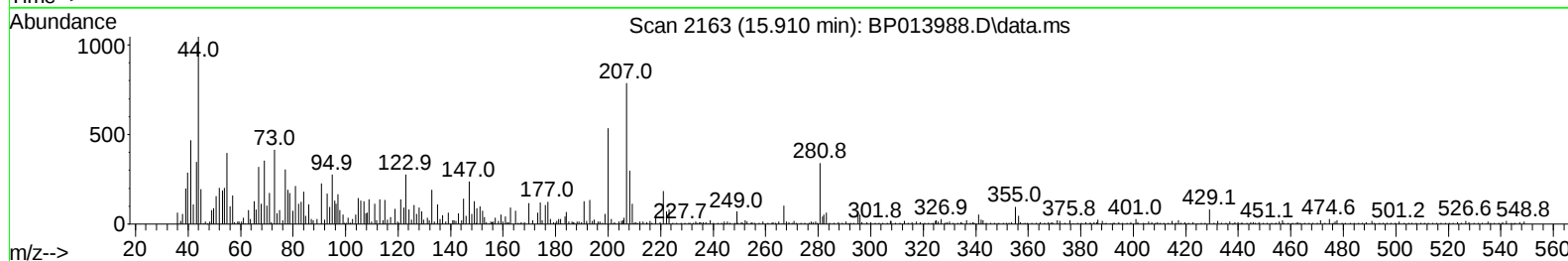
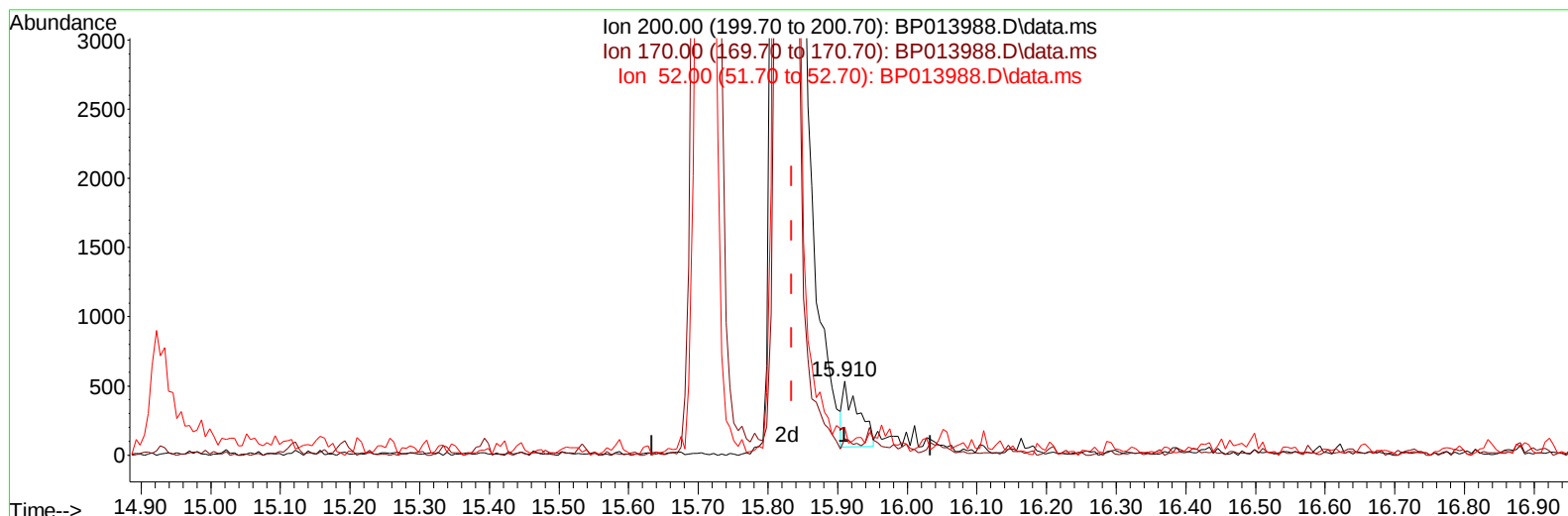
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TIC: BP013988.D\data.ms

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

15.910min (+ 0.076) 0.14 ng/u1

response 706

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.80	21.72
52.00	37.50	37.64
0.00	0.00	0.00

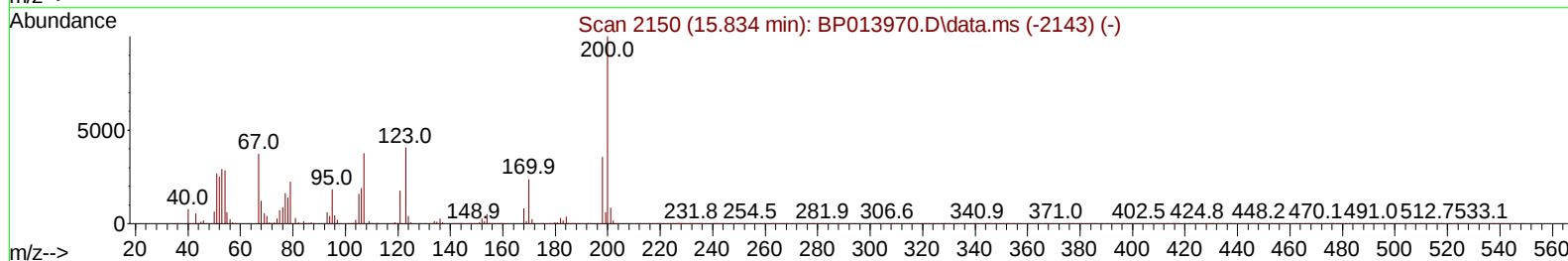
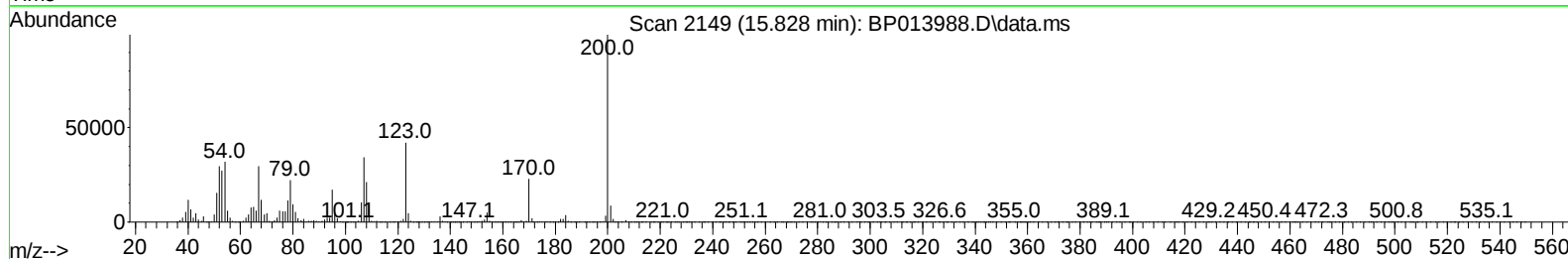
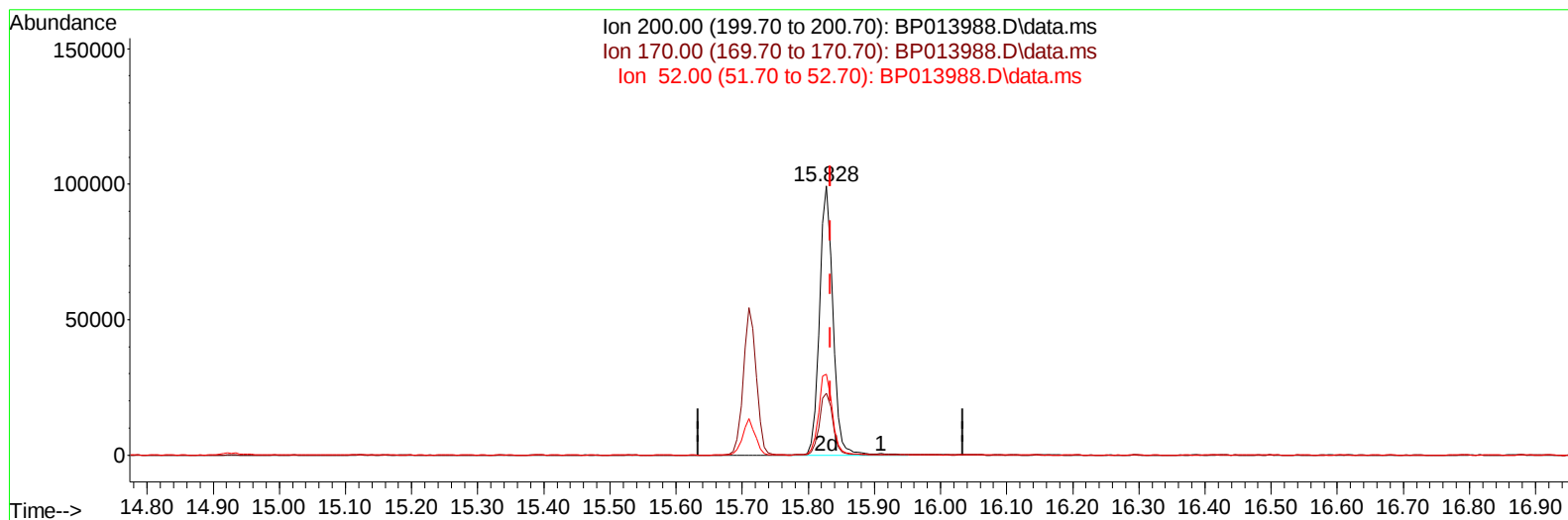
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TIC: BP013988.D\data.ms

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

15.828min (-0.006) 27.28 ng/ul m

response 138391

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.80	23.05
52.00	37.50	29.96#
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	86505	20.000	ng/ul	0.00
20) Naphthalene-d8	10.904	136	361932	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.722	164	265565	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.475	188	671312	20.000	ng/ul	0.00
79) Chrysene-d12	21.551	240	802785	20.000	ng/ul	0.00
88) Perylene-d12	24.086	264	904193	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	10598	4.629	ng/uL	0.00
4) Pyridine-d5	3.875	84	56786	8.958	ng/ul	0.00
7) Phenol-d5	7.234	99	38943	5.101	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.399	67	133636	30.017	ng/ul	0.00
11) 2-Chlorophenol-d4	7.610	132	125523	21.582	ng/ul	0.00
15) 4-Methylphenol-d8	8.793	113	83465	13.385	ng/ul	0.00
21) Nitrobenzene-d5	9.257	128	85338	29.551	ng/ul	0.00
24) 2-Nitrophenol-d4	9.981	143	94900	27.452	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.522	165	169792	26.104	ng/ul	0.00
31) 4-Chloroaniline-d4	11.046	131	217975	25.323	ng/ul	0.00
46) Dimethylphthalate-d6	14.122	166	726500	32.235	ng/ul	0.00
49) Acenaphthylene-d8	14.416	160	729293	30.546	ng/ul	0.00
54) 4-Nitrophenol-d4	14.922	143	16457m	4.144	ng/ul	0.00
60) Fluorene-d10	15.710	176	612694	32.077	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.828	200	138391m	27.276	ng/ul	0.00
73) Anthracene-d10	17.575	188	1087672	33.375	ng/ul	0.00
81) Pyrene-d10	19.792	212	1446607	32.672	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.921	264	1625586	33.799	ng/ul	0.00

Target Compounds Qvalue

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

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