

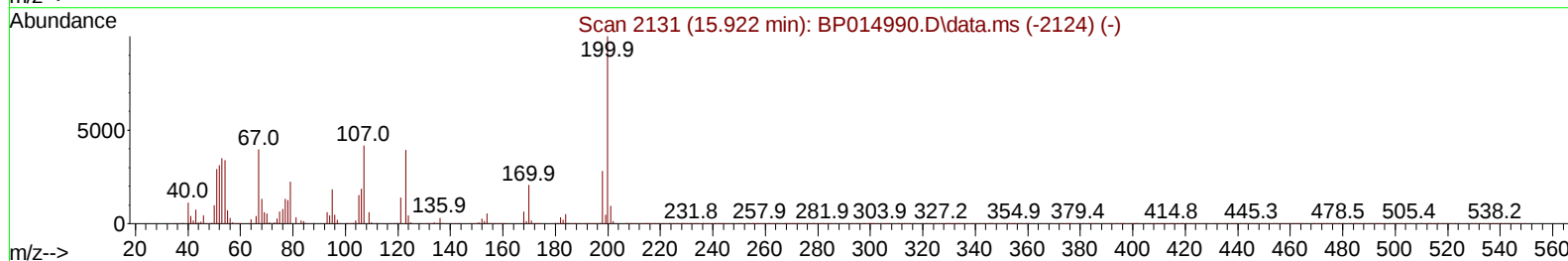
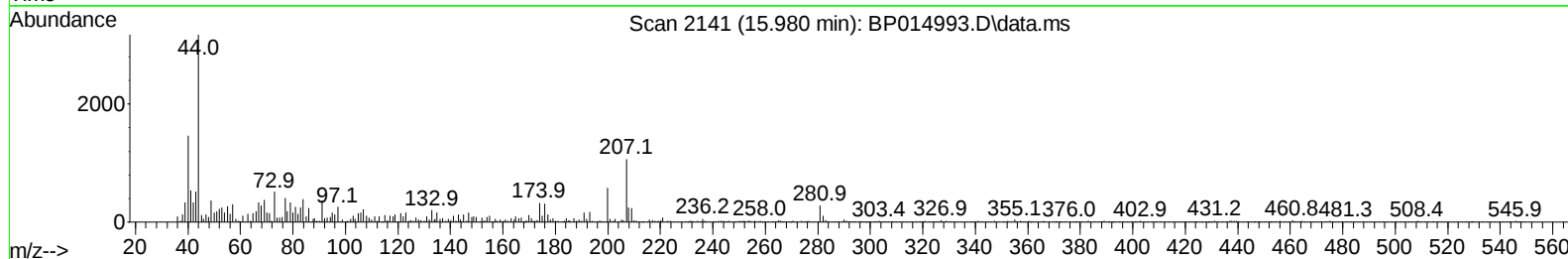
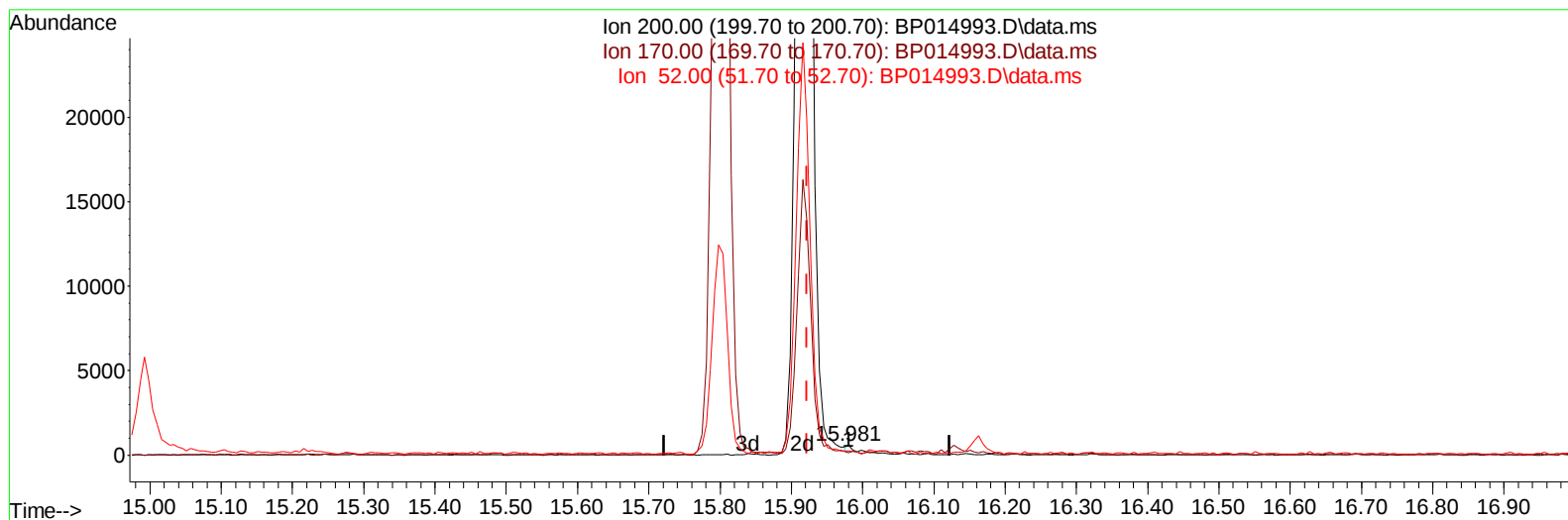
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP050823\
 Data File : BP014993.D
 Acq On : 08 May 2023 13:43
 Operator : CG/JU
 Sample : 02589-23
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 JRED5

Manual IntegrationsAPPROVED

Quant Time: May 08 22:50:43 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP050323.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri May 05 12:12:36 2023
 Response via : Initial Calibration

Reviewed By :Christian Giraldo 05/09/2023
 Supervised By :Jagrut Upadhyay 05/09/2023



TIC: BP014993.D\data.ms

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

15.980min (+ 0.059) 0.02 ng/u1

response 169

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.30	20.24
52.00	44.00	39.79
0.00	0.00	0.00

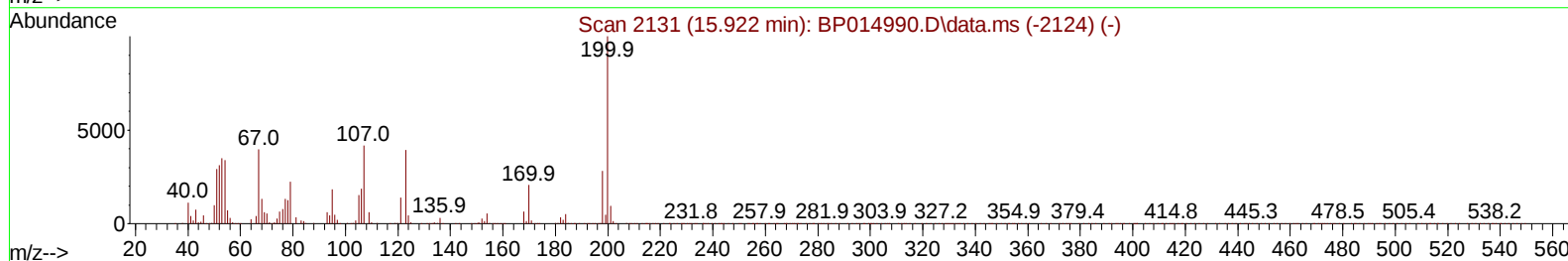
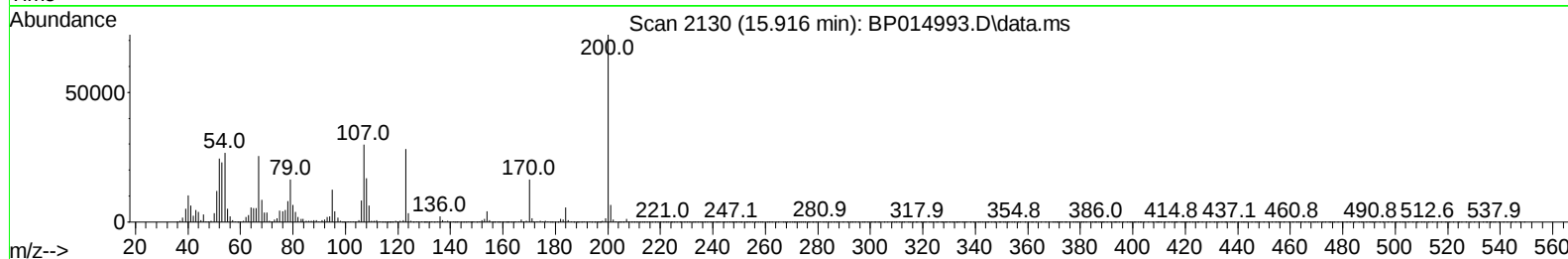
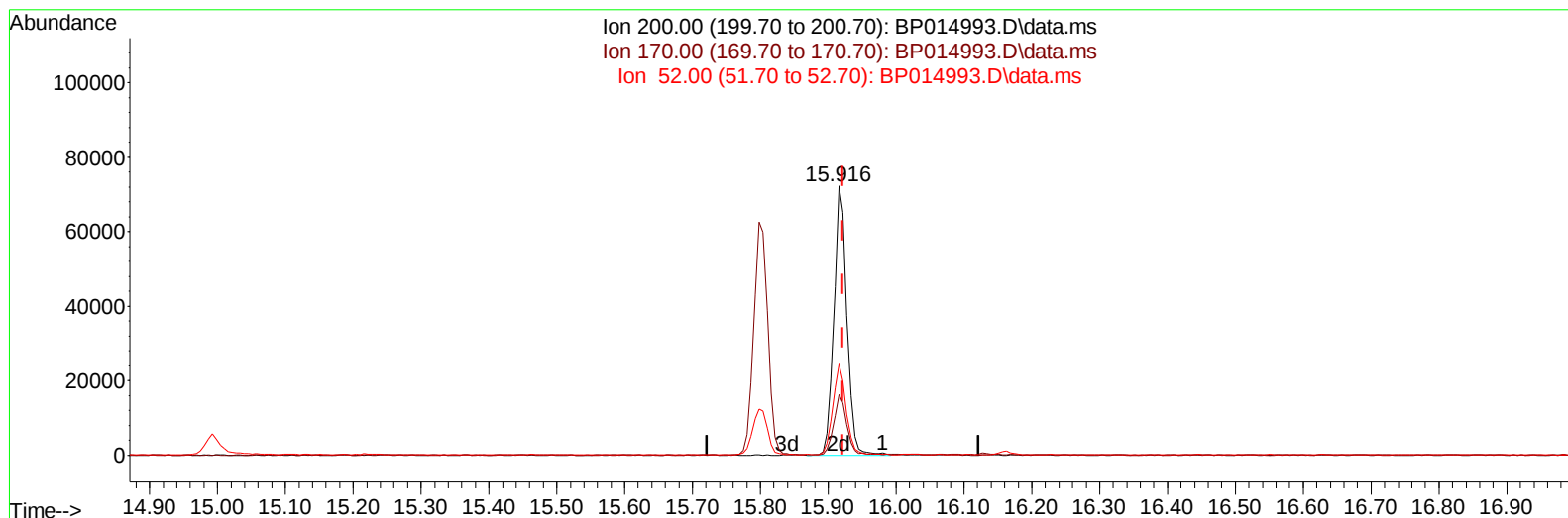
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TIC: BP014993.D\data.ms

(65) 4,6-Dinitro-2-methylphenol-d2 (S)**15.916min (-0.006) 8.98 ng/ul m****response 96902**

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	19.30	22.58
52.00	44.00	33.83#
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	8.163	152	306135	20.000	ng/ul	0.00
20) Naphthalene-d8	10.998	136	1398154	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.810	164	907143	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.563	188	1916206	20.000	ng/ul	0.00
79) Chrysene-d12	21.645	240	1269259	20.000	ng/ul	0.00
88) Perylene-d12	24.251	264	1131551	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.458	96	14546	1.767	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.310	99	318763	12.010	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.481	67	203095	12.380	ng/ul	0.00
11) 2-Chlorophenol-d4	7.687	132	252617	12.944	ng/ul	0.00
15) 4-Methylphenol-d8	8.875	113	242688	11.865	ng/ul	0.00
21) Nitrobenzene-d5	9.340	128	130407	13.114	ng/ul	0.00
24) 2-Nitrophenol-d4	10.069	143	143872	14.090	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.610	165	261627	13.049	ng/ul	0.00
31) 4-Chloroaniline-d4	11.134	131	305024	10.518	ng/ul	0.00
46) Dimethylphthalate-d6	14.210	166	860199	13.398	ng/ul	0.00
49) Acenaphthylene-d8	14.504	160	1014654	13.177	ng/ul	0.00
54) 4-Nitrophenol-d4	14.992	143	74917	6.707	ng/ul	0.00
60) Fluorene-d10	15.798	176	744509	13.479	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.916	200	96902m	8.983	ng/ul	0.00
73) Anthracene-d10	17.663	188	1131450	13.141	ng/ul	0.00
81) Pyrene-d10	19.880	212	1227455	14.339	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.074	264	754279	13.681	ng/ul	0.00

Target Compounds Qvalue

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

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