

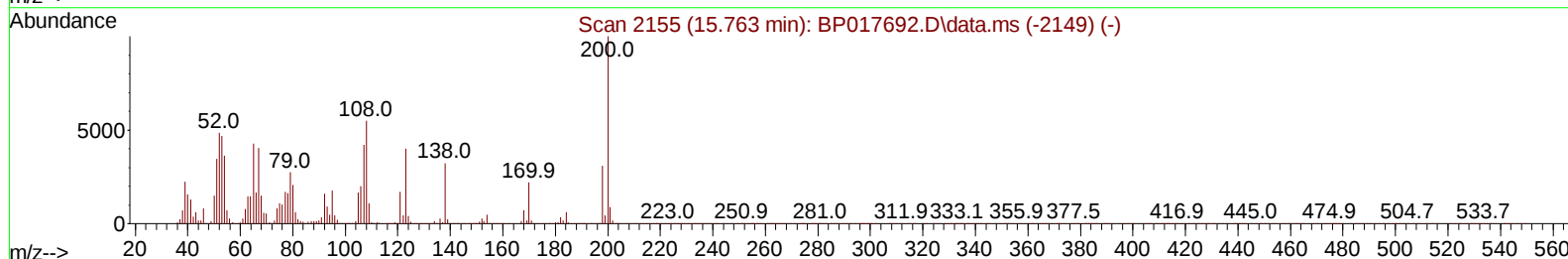
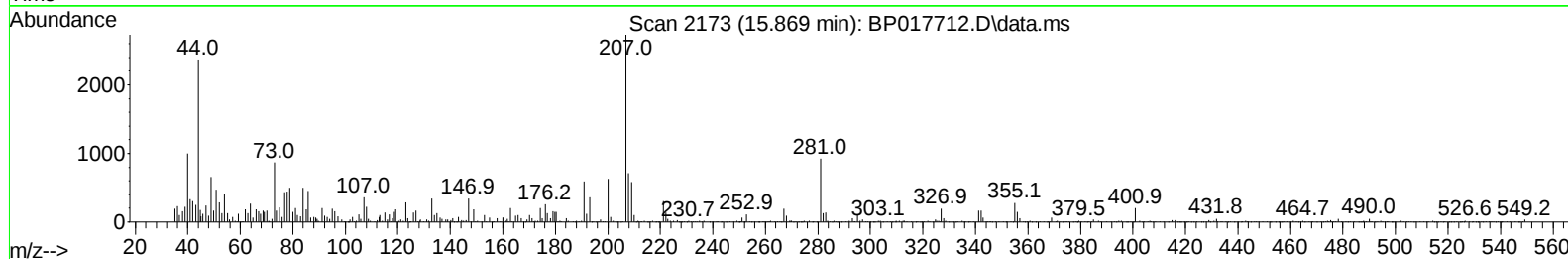
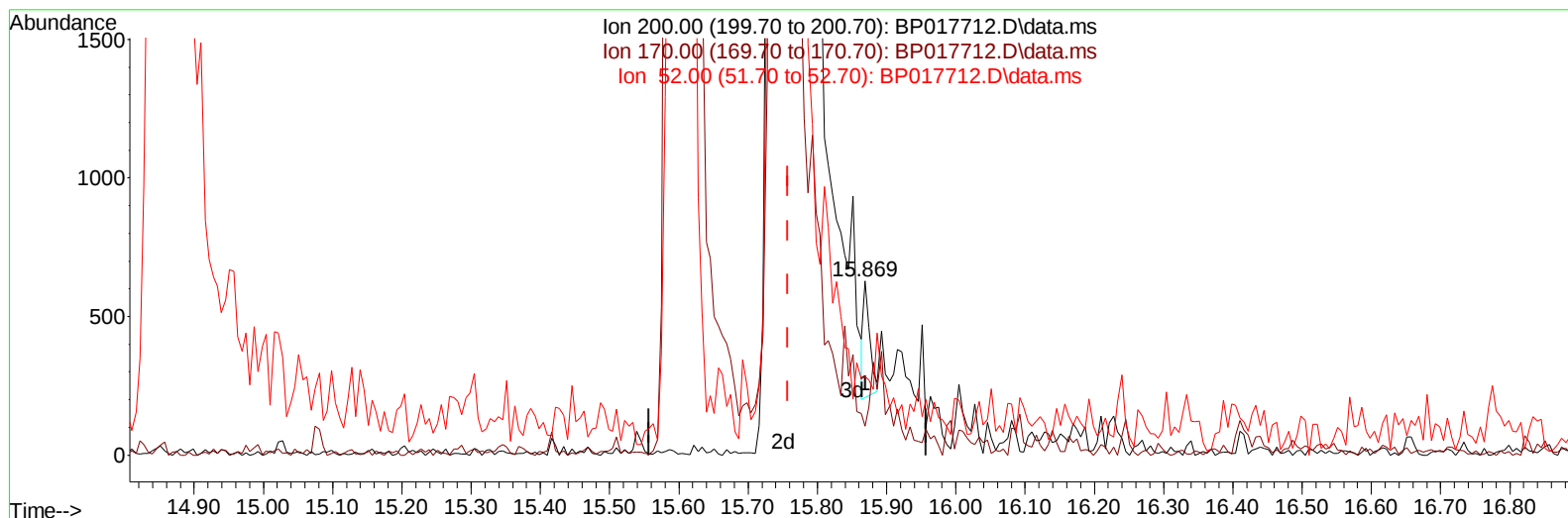
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP101723\
 Data File : BP017712.D
 Acq On : 17 Oct 2023 23:14
 Operator : MA/JU
 Sample : PB156379BL
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK379

Manual IntegrationsAPPROVED

Quant Time: Oct 18 00:46:57 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP101223.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Oct 12 16:14:36 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/18/2023
 Supervised By :mohammad ahmed 10/19/2023



TIC: BP017712.D\data.ms

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

15.869min (+ 0.112) 0.06 ng/u1

response 279

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	20.20	16.56
52.00	42.70	45.06
0.00	0.00	0.00

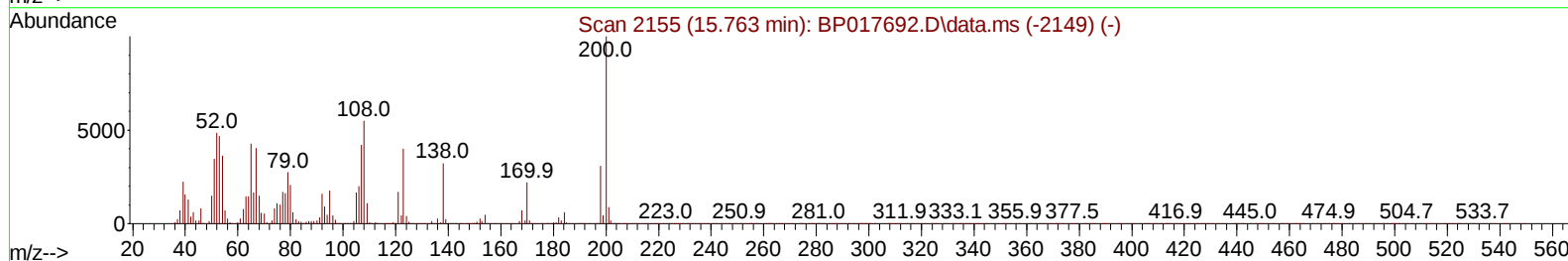
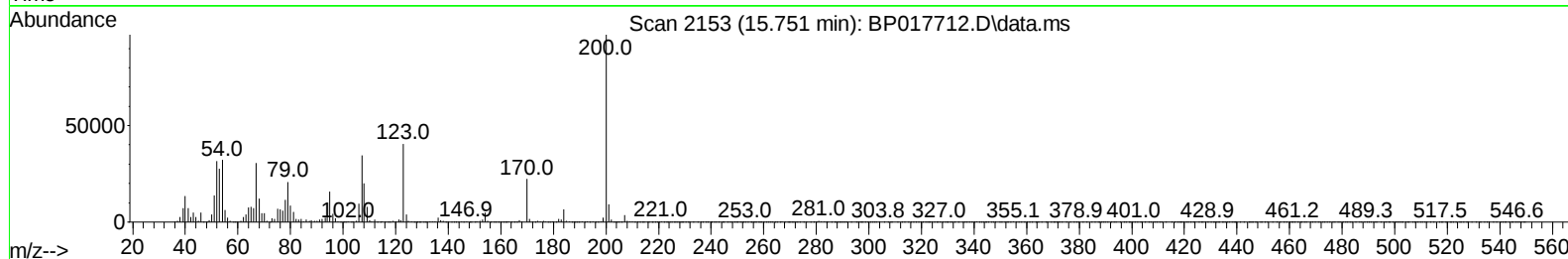
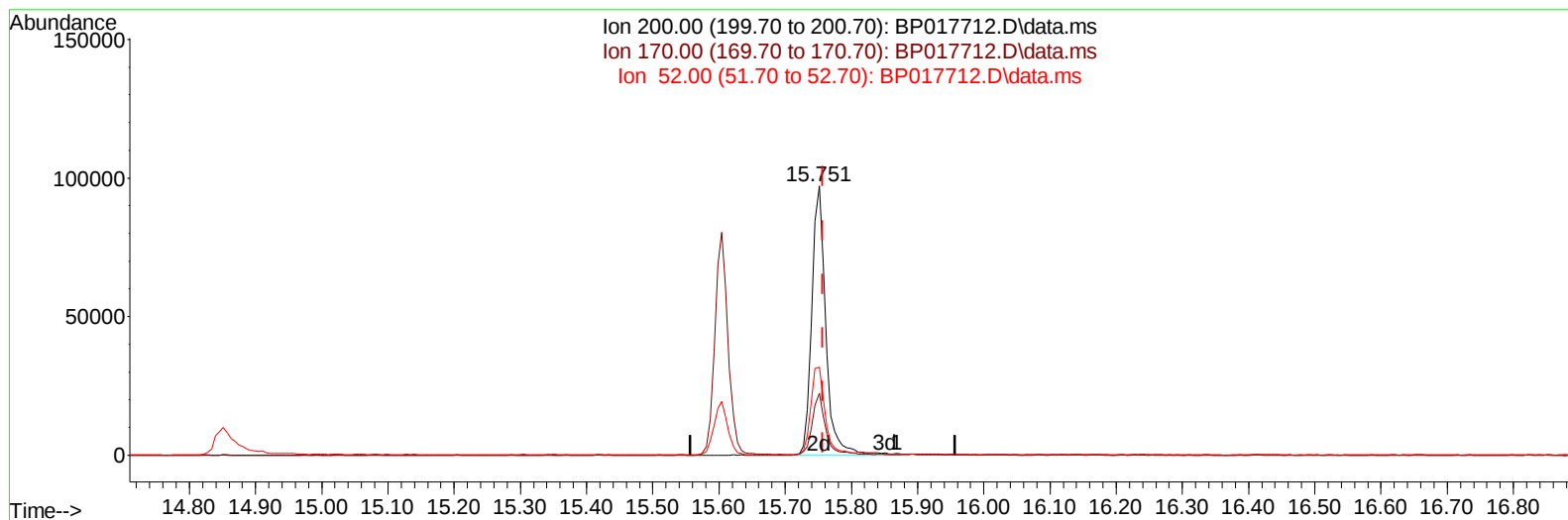
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.751min (-0.006) 29.81 ng/ul m

response 141623

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	20.20	22.99
52.00	42.70	32.64#
0.00	0.00	0.00

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Quant Time: Oct 18 01:18:39 2023
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.899	152	114632	20.000	ng/ul	0.00
20) Naphthalene-d8	10.734	136	463985	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.598	164	289689	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.392	188	672067	20.000	ng/ul	0.00
79) Chrysene-d12	21.533	240	715132	20.000	ng/ul	-0.01
88) Perylene-d12	24.098	264	841114	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.258	96	24694	7.942	ng/uL	-0.01
4) Pyridine-d5	3.693	84	338595	39.395	ng/ul	0.00
7) Phenol-d5	7.052	99	436026	40.856	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.240	67	273790	42.303	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.416	132	354859	42.957	ng/ul	0.00
15) 4-Methylphenol-d8	8.622	113	342851	40.220	ng/ul	0.00
21) Nitrobenzene-d5	9.104	128	164336	42.168	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.822	143	183302	41.824	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.351	165	337069	40.921	ng/ul	0.00
31) 4-Chloroaniline-d4	10.904	131	475323	40.259	ng/ul	0.00
46) Dimethylphthalate-d6	14.016	166	1069083	42.611	ng/ul	0.00
49) Acenaphthylene-d8	14.298	160	1165420	43.136	ng/ul	0.00
54) 4-Nitrophenol-d4	14.851	143	153545	36.593	ng/ul	0.00
60) Fluorene-d10	15.604	176	938829	43.990	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.751	200	141623m	29.814	ng/ul	0.00
73) Anthracene-d10	17.492	188	1519325	43.791	ng/ul	0.00
81) Pyrene-d10	19.751	212	1900386	44.153	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.927	264	2081428	44.080	ng/ul	-0.01

Target Compounds Qvalue

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

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