

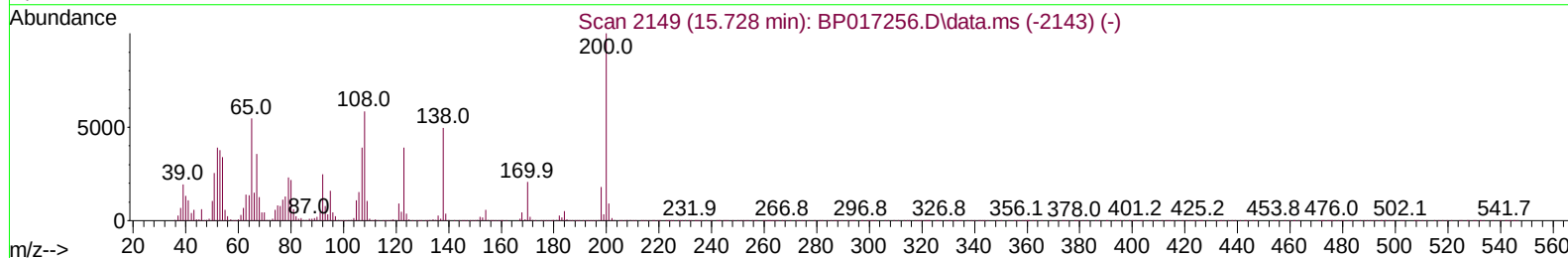
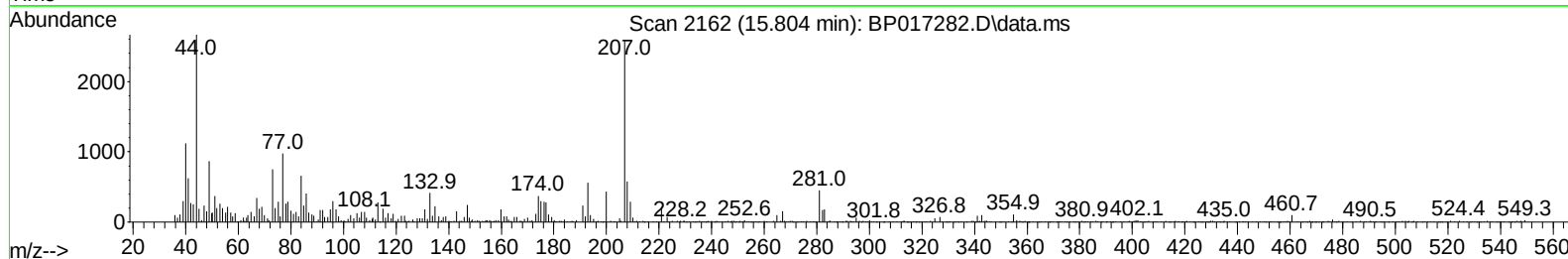
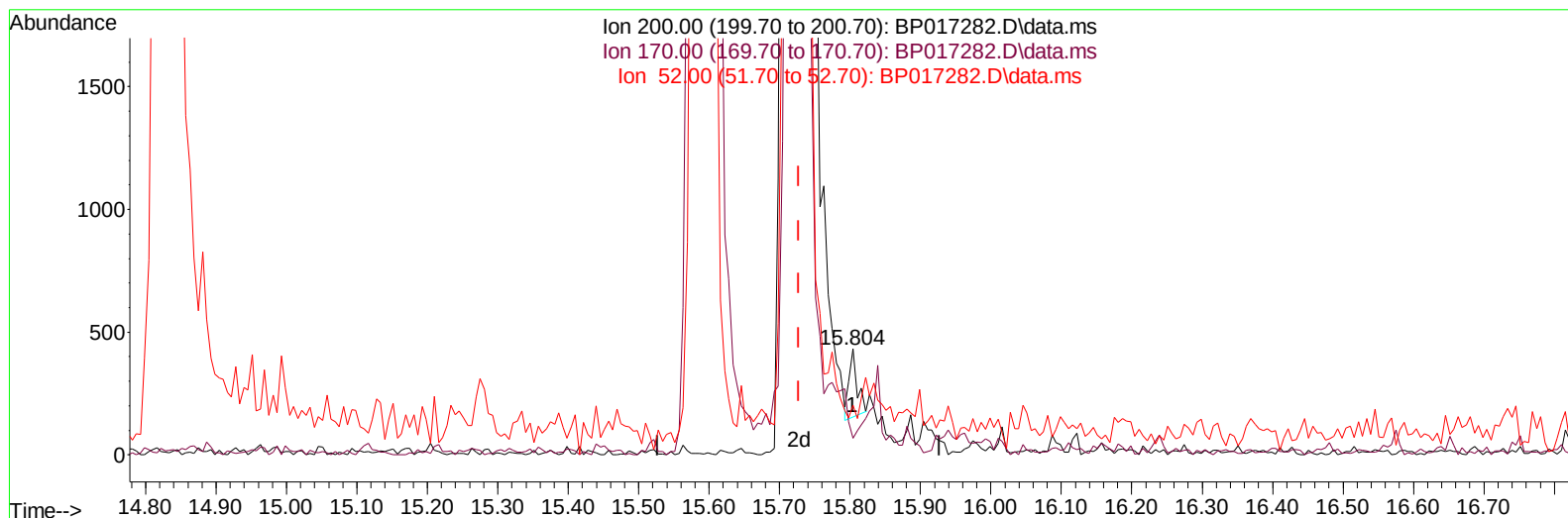
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP092223\
 Data File : BP017282.D
 Acq On : 22 Sep 2023 10:04
 Operator : MA/JU
 Sample : PB155610BL
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK610

Manual IntegrationsAPPROVED

Quant Time: Sep 22 10:45:15 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Sep 21 23:51:07 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 09/25/2023
 Supervised By :mohammad ahmed 09/27/2023



TIC: BP017282.D\data.ms

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

15.804min (+ 0.077) 0.05 ng/ul

response 217

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	18.80	15.55
52.00	41.50	47.10
0.00	0.00	0.00

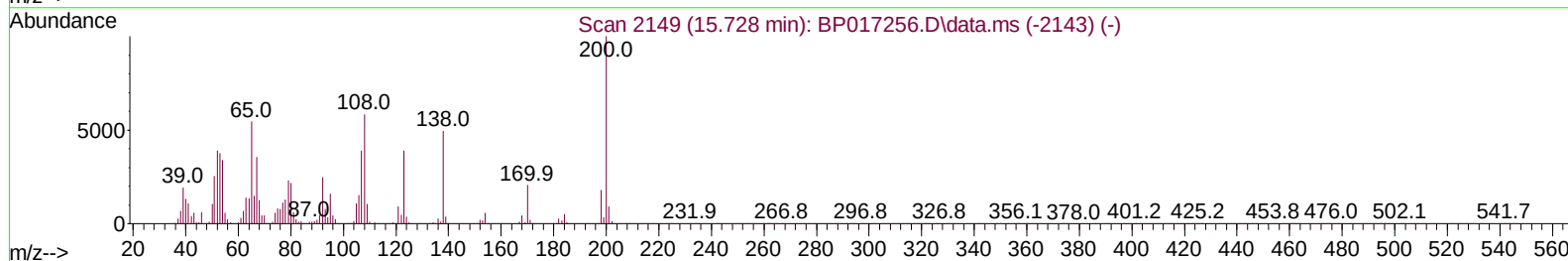
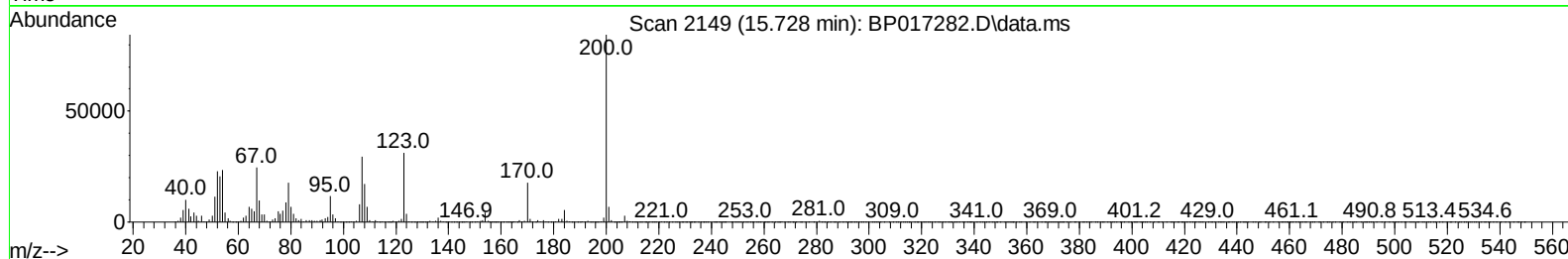
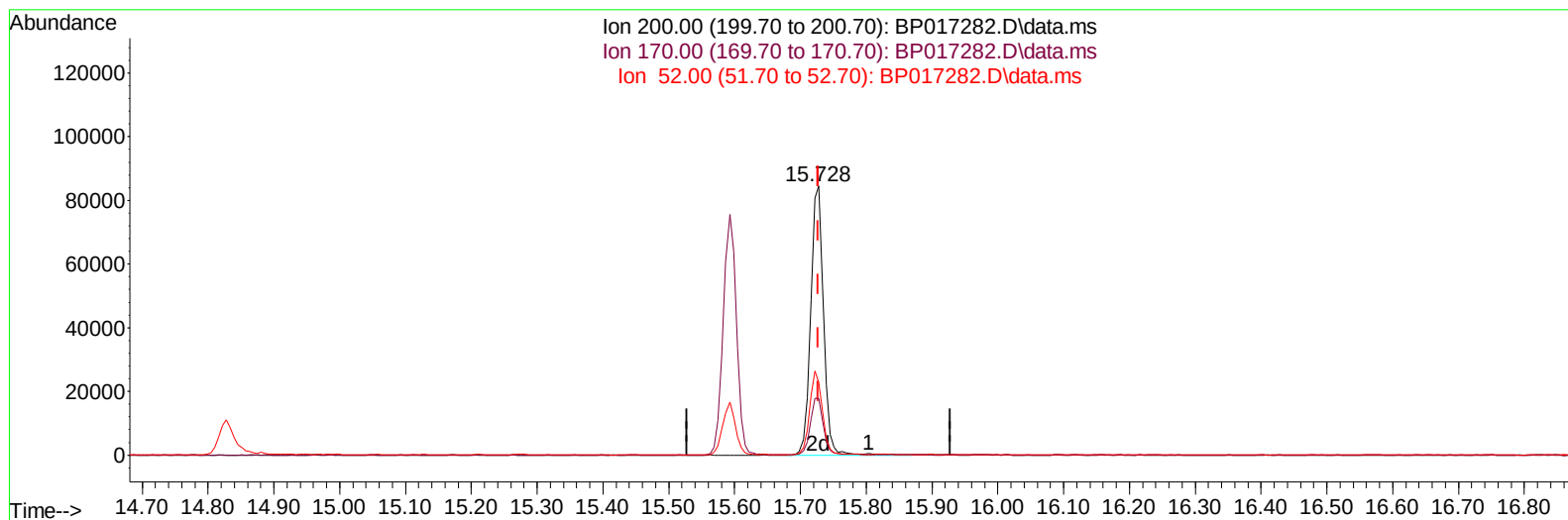
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TIC: BP017282.D\data.ms

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

15.728min (+ 0.000) 26.34 ng/ul m

response 116012

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	18.80	21.03
52.00	41.50	27.06#
0.00	0.00	0.00

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Quant Time: Sep 22 11:05:22 2023
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.934	152		151514	20.000	ng/ul	0.00
20) Naphthalene-d8	10.757	136		599597	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.598	164		345129	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.369	188		781434	20.000	ng/ul	0.00
79) Chrysene-d12	21.469	240		790252	20.000	ng/ul	0.00
88) Perylene-d12	23.992	264		893490	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.305	96		28522	7.732	ng/uL	0.00
4) Pyridine-d5	3.740	84		394437	38.249	ng/ul	0.00
7) Phenol-d5	7.093	99		467763	37.590	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.275	67		280060	37.293	ng/ul	0.00
11) 2-Chlorophenol-d4	7.458	132		379963	38.498	ng/ul	0.00
15) 4-Methylphenol-d8	8.652	113		358683	37.075	ng/ul	0.00
21) Nitrobenzene-d5	9.134	128		173781	40.090	ng/ul	0.00
24) 2-Nitrophenol-d4	9.846	143		190846	40.548	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.375	165		342331	38.497	ng/ul	0.00
31) 4-Chloroaniline-d4	10.922	131		484984	37.650	ng/ul	0.00
46) Dimethylphthalate-d6	14.010	166		996208	40.293	ng/ul	0.00
49) Acenaphthylene-d8	14.298	160		1093606	38.459	ng/ul	0.00
54) 4-Nitrophenol-d4	14.828	143		156563	36.348	ng/ul	0.00
60) Fluorene-d10	15.592	176		862387	40.516	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.728	200		116012m	26.340	ng/ul	0.00
73) Anthracene-d10	17.469	188		1459708	42.119	ng/ul	0.00
81) Pyrene-d10	19.710	212		1806835	42.280	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.821	264		1843550	42.854	ng/ul	-0.01

Target Compounds Qvalue

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

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