

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
 Data File : BP018229.D
 Acq On : 17 Nov 2023 20:15
 Operator : MA/JU
 Sample : PB157215BL
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

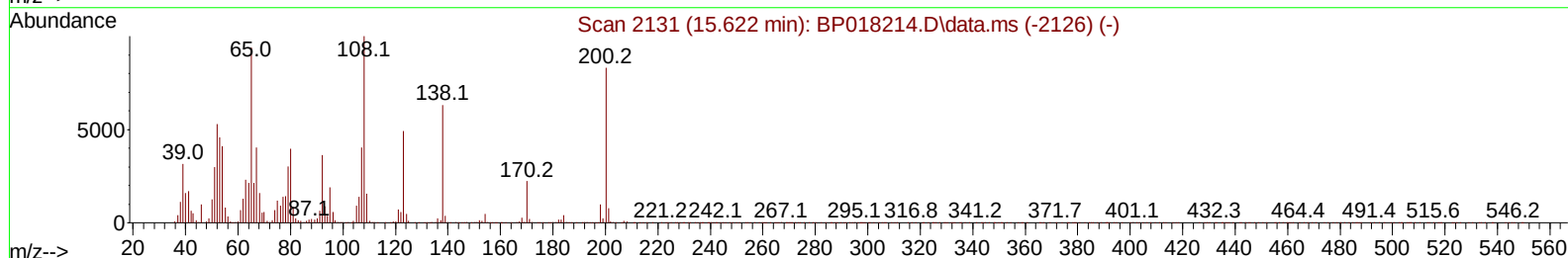
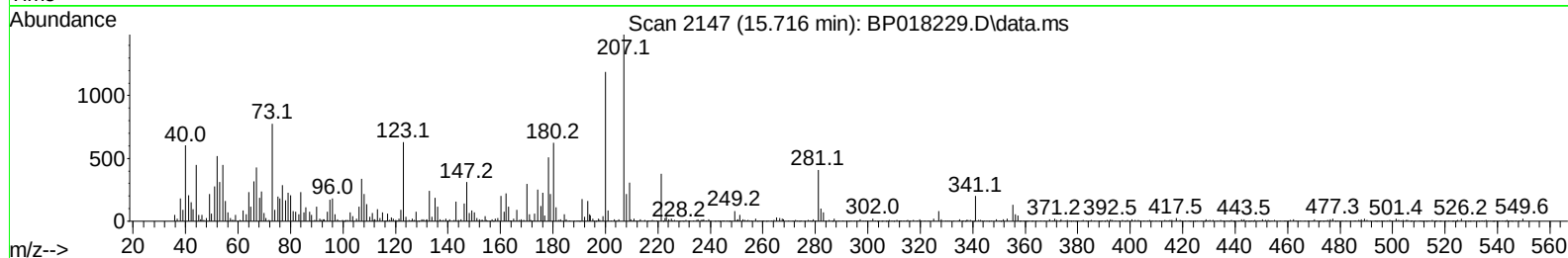
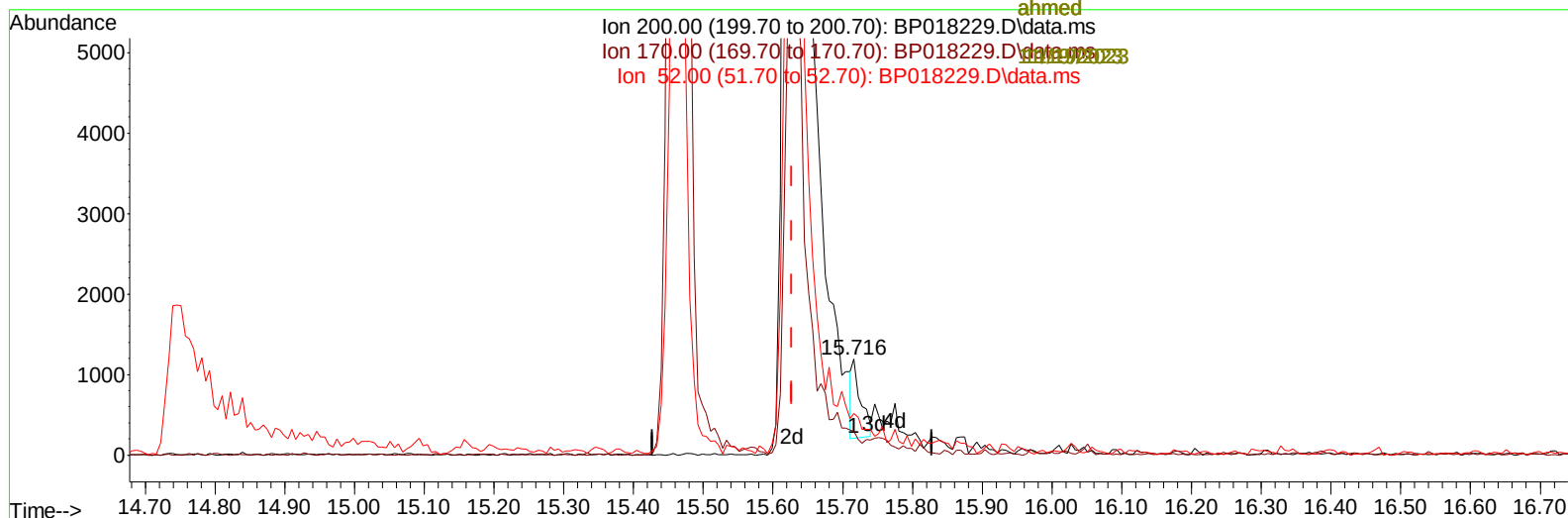
Instrument :
 BNA_P
 ClientSampleId :
 SBLK215

Manual IntegrationsAPPROVED

Quant Time: Nov 17 21:53:48 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
 Response via : Initial Calibration

Reviewed By :Yogesh
 Patel

11/19/2023
 Supervised By :mohammad
 ahmed



TIC: BP018229.D\data.ms

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

15.716min (+ 0.088) 0.58 ng/ul

response 845

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	24.87
52.00	52.00	43.53
0.00	0.00	0.00

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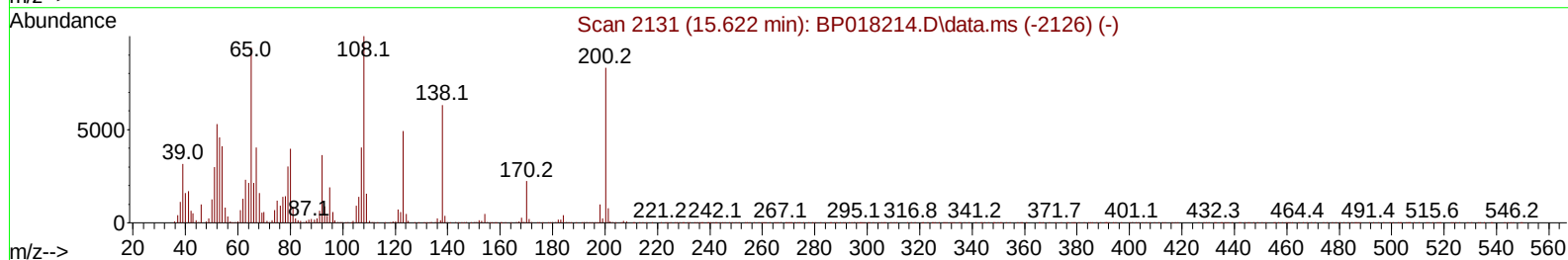
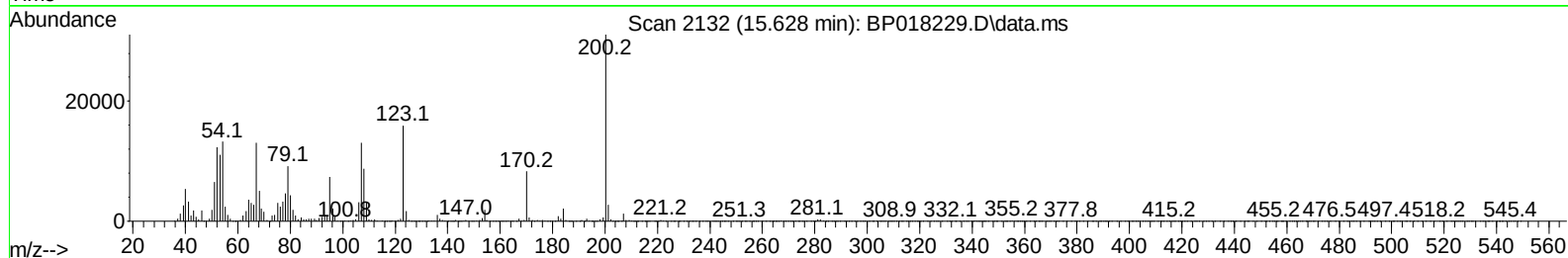
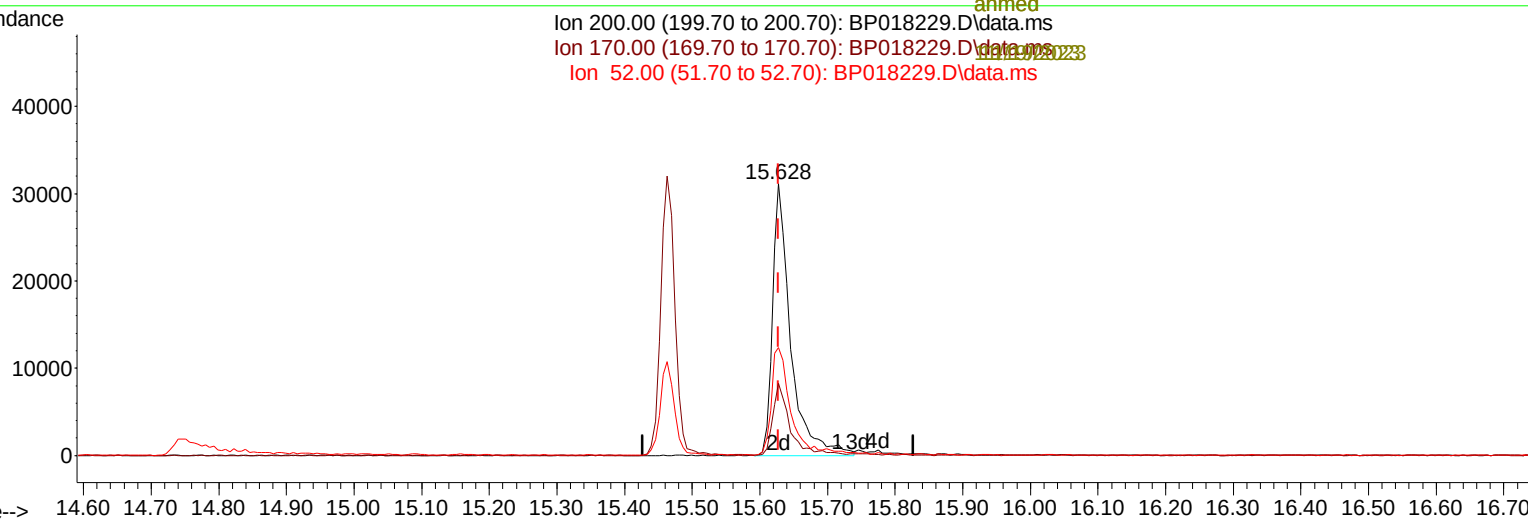
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Ion 200.00 (199.70 to 200.70): BP018229.D\data.ms
 Ion 170.00 (169.70 to 170.70): BP018229.D\data.ms
 Ion 52.00 (51.70 to 52.70): BP018229.D\data.ms



TIC: BP018229.D\data.ms

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

15.628min (+ 0.000) 39.48 ng/ul m

response 57436

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	26.66
52.00	52.00	39.79#
0.00	0.00	0.00

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Quant Time: Nov 17 22:32:27 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 15 18:13:37 2023
 Response via : Initial Calibration

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----11/19/2023							
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.781	152		42081	20.000	ng/ul	0.00
20) Naphthalene-d8	10.599	136		148309	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.463	164		90020	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.245	188		201260	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240		179550	20.000	ng/ul	0.00
88) Perylene-d12	23.851	264		191493	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.158	96		11602	9.726	ng/uL	0.00
4) Pyridine-d5	3.593	84		129194	41.537	ng/ul	0.00
7) Phenol-d5	6.952	99		152677	37.581	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67		97716	40.928	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132		133212	41.406	ng/ul	0.00
15) 4-Methylphenol-d8	8.505	113		130147	39.094	ng/ul	0.00
21) Nitrobenzene-d5	8.987	128		63118	47.238	ng/ul	0.00
24) 2-Nitrophenol-d4	9.705	143		71844	47.460	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165		114231	42.339	ng/ul	0.00
31) 4-Chloroaniline-d4	10.781	131		156607	42.327	ng/ul	0.00
46) Dimethylphthalate-d6	13.887	166		441690	53.178	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160		451406	49.898	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143		35817	28.858	ng/ul	0.01
60) Fluorene-d10	15.463	176		358431	52.269	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200		57436m	39.476	ng/ul	0.00
73) Anthracene-d10	17.345	188		570661	52.427	ng/ul	0.00
81) Pyrene-d10	19.598	212		684646	57.086	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.680	264		626344	56.251	ng/ul	0.00

Target Compounds Qvalue

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

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