

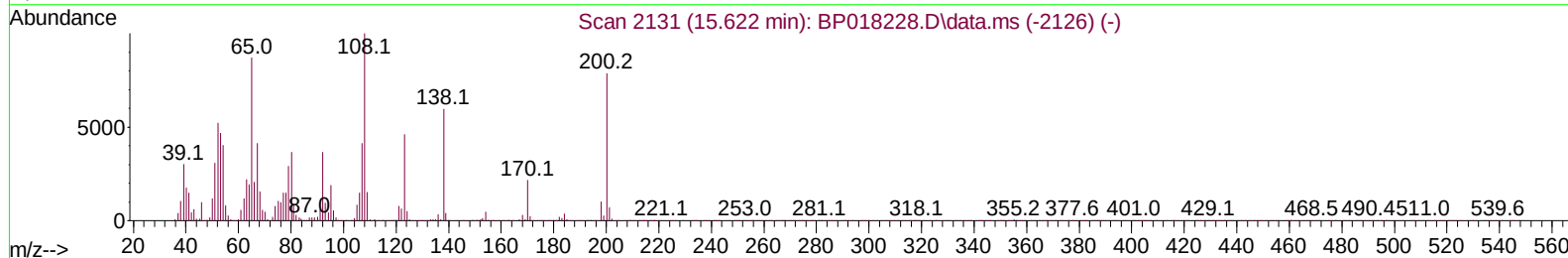
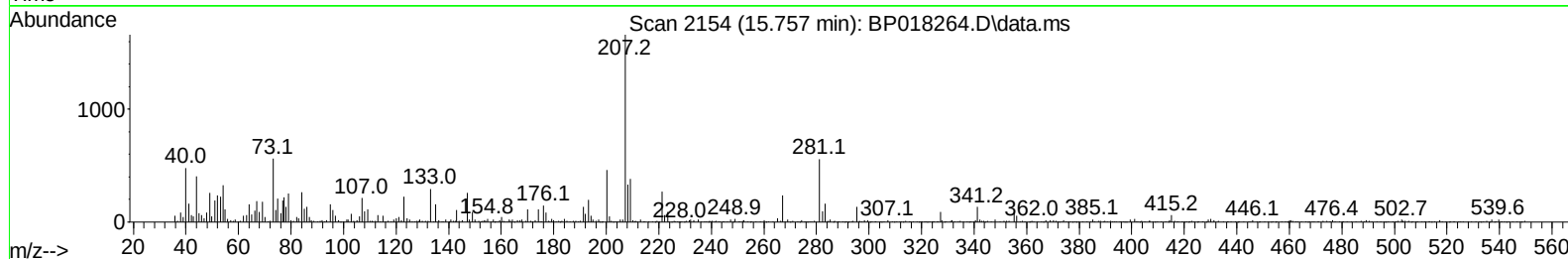
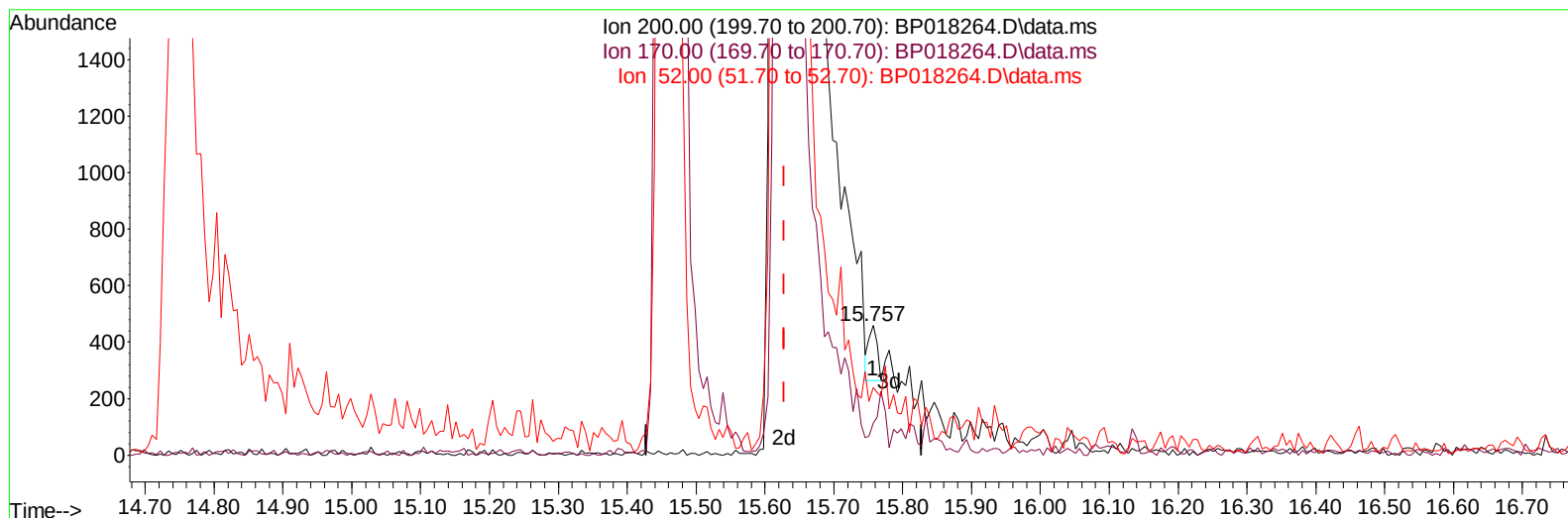
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111723\
Data File : BP018264.D
Acq On : 18 Nov 2023 16:38
Operator : MA/JU
Sample : PB157225BL
Misc :
ALS Vial : 53 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK225

Manual IntegrationsAPPROVED

Quant Time: Nov 19 00:56:15 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 11/19/2023



TIC: BP018264.D\data.ms

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

15.757min (+ 0.129) 0.07 ng/u1

response 167

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	24.84
52.00	52.00	52.07
0.00	0.00	0.00

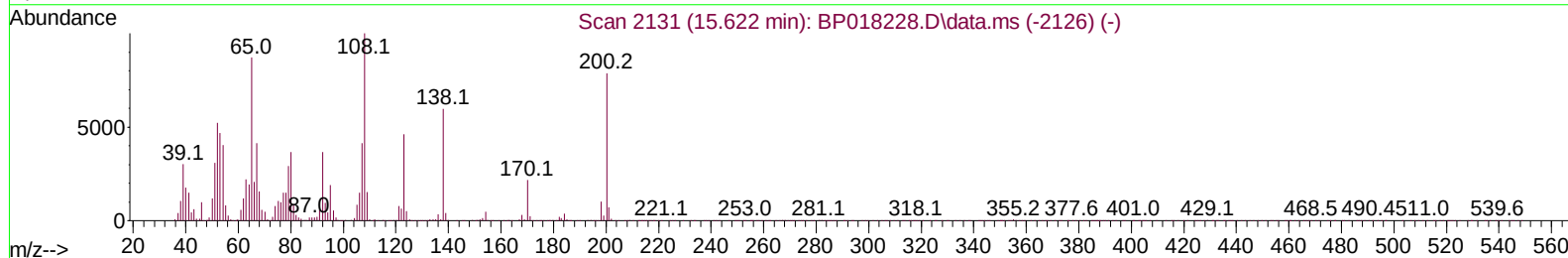
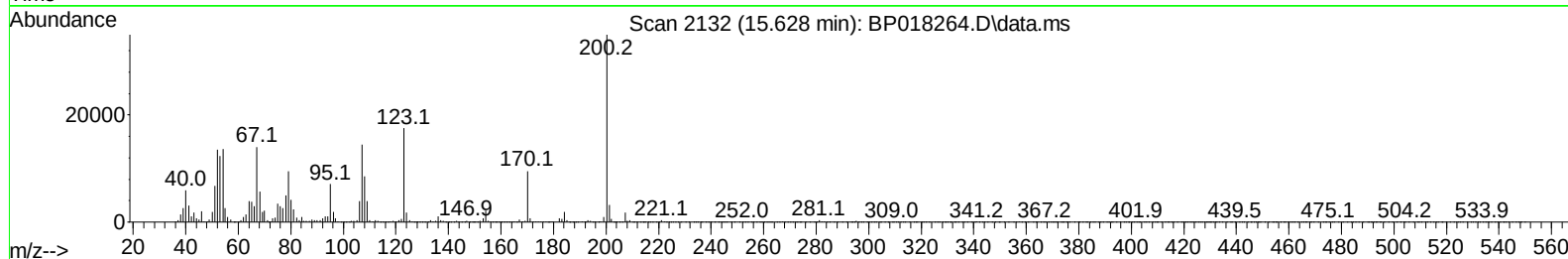
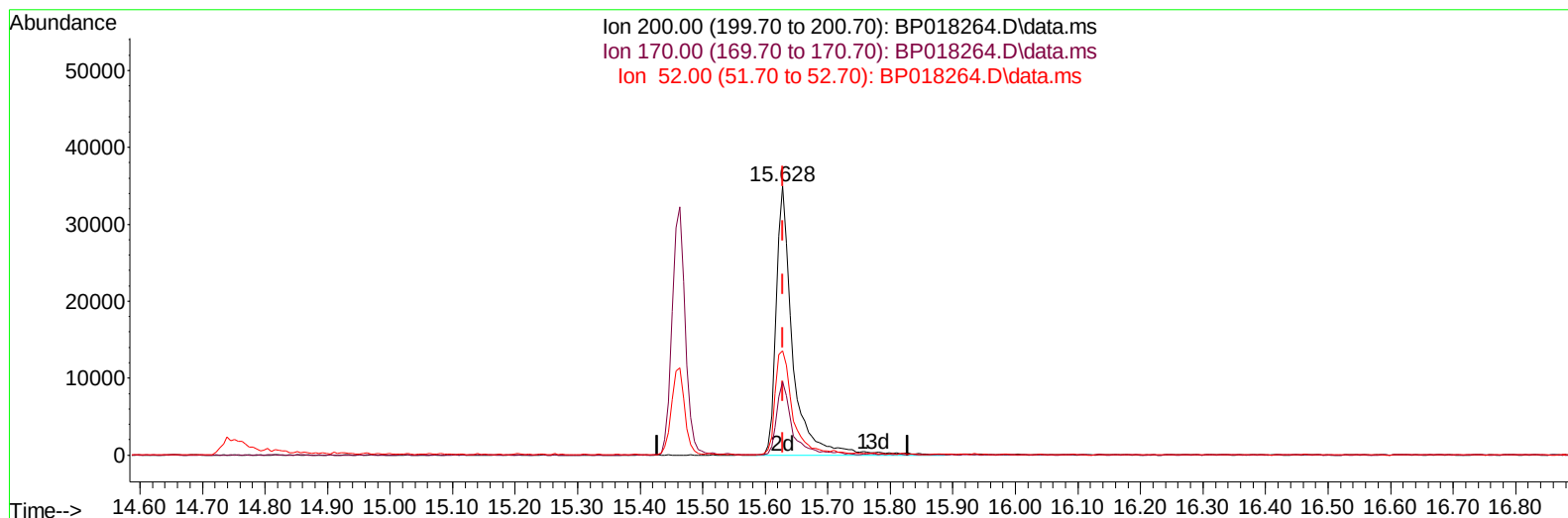
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TIC: BP018264.D\data.ms

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

15.628min (+ 0.000) 25.80 ng/ul m

response 64221

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	27.28
52.00	52.00	38.76#
0.00	0.00	0.00

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Instrument :
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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/19/2023
 Supervised By :mohammad ahmed 11/19/2023

Quant Time: Nov 19 02:48:53 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
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.775	152	67171	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.593	136	253656	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	153305	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.239	188	344300	20.000	ng/ul	0.00
79) Chrysene-d12	21.380	240	335408	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	367531	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.158	96	11786	6.190	ng/uL	0.00
4) Pyridine-d5	3.593	84	131847	26.556	ng/ul	0.00
7) Phenol-d5	6.946	99	165020	25.447	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.128	67	102033	26.773	ng/ul	0.00
11) 2-Chlorophenol-d4	7.299	132	143879	28.017	ng/ul	0.00
15) 4-Methylphenol-d8	8.505	113	130744	24.604	ng/ul	0.00
21) Nitrobenzene-d5	8.981	128	63679	27.865	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.699	143	69975	27.027	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.228	165	118930	25.773	ng/ul	0.00
31) 4-Chloroaniline-d4	10.781	131	169891	26.847	ng/ul	0.00
46) Dimethylphthalate-d6	13.881	166	434481	30.716	ng/ul	0.00
49) Acenaphthylene-d8	14.157	160	461363	29.946	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	39010	18.456	ng/ul	0.01
60) Fluorene-d10	15.463	176	361749	30.976	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.628	200	64221m	25.802	ng/ul	0.00
73) Anthracene-d10	17.339	188	576176	30.942	ng/ul	0.00
81) Pyrene-d10	19.598	212	726019	32.406	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.686	264	713527	33.388	ng/ul	0.00

Target Compounds Qvalue

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

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