

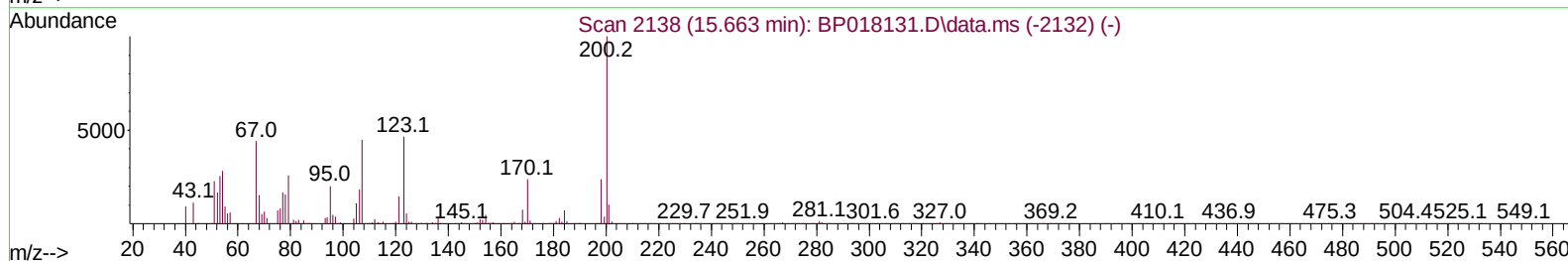
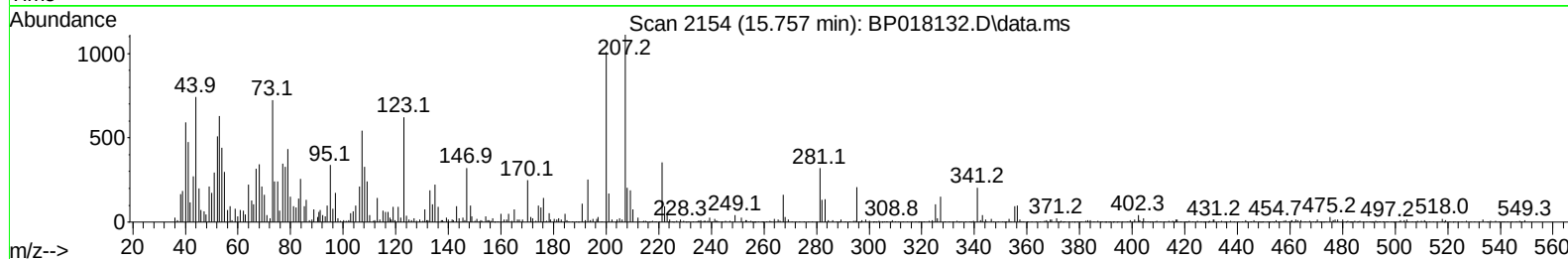
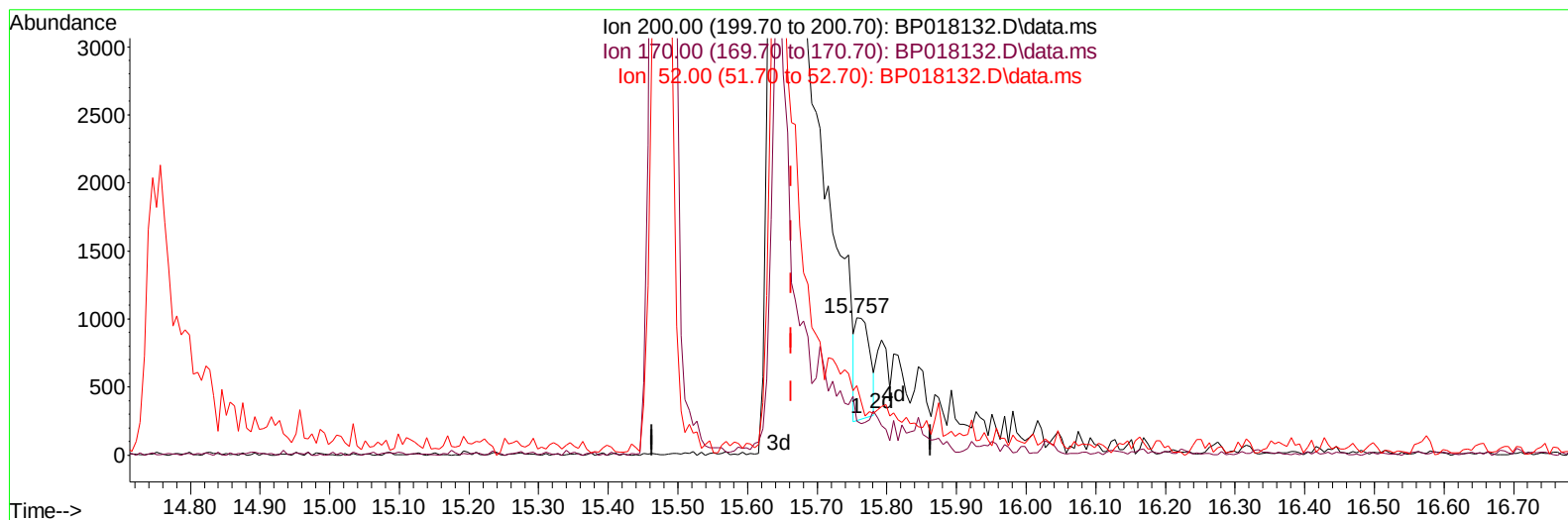
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111323\
 Data File : BP018132.D
 Acq On : 13 Nov 2023 15:44
 Operator : MA/JU
 Sample : PB157055BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK055

Manual IntegrationsAPPROVED

Quant Time: Nov 13 17:42:28 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 13 15:50:18 2023
 Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/14/2023
 Supervised By :mohammad ahmed 11/16/2023



TIC: BP018132.D\data.ms

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

15.757min (+ 0.094) 0.35 ng/ul

response 1068

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	24.58
52.00	52.00	50.35
0.00	0.00	0.00

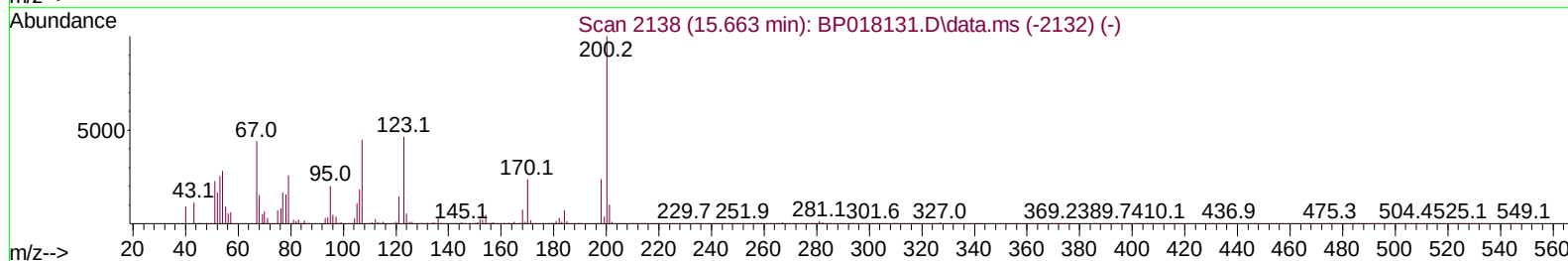
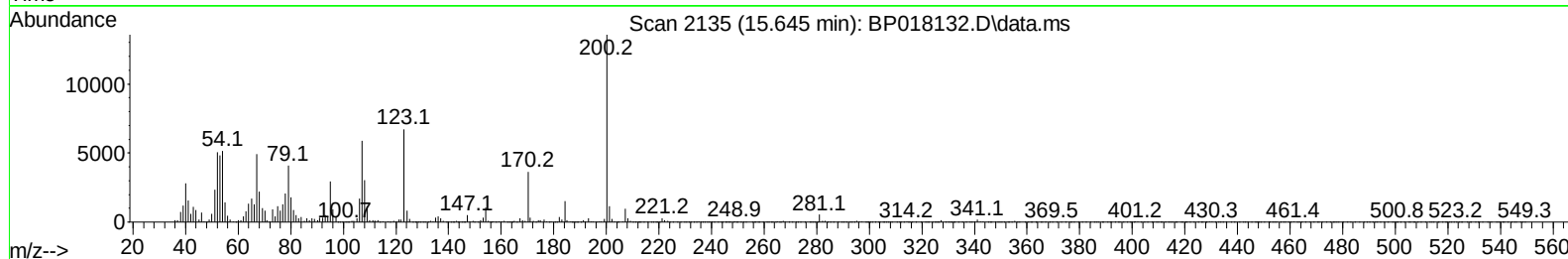
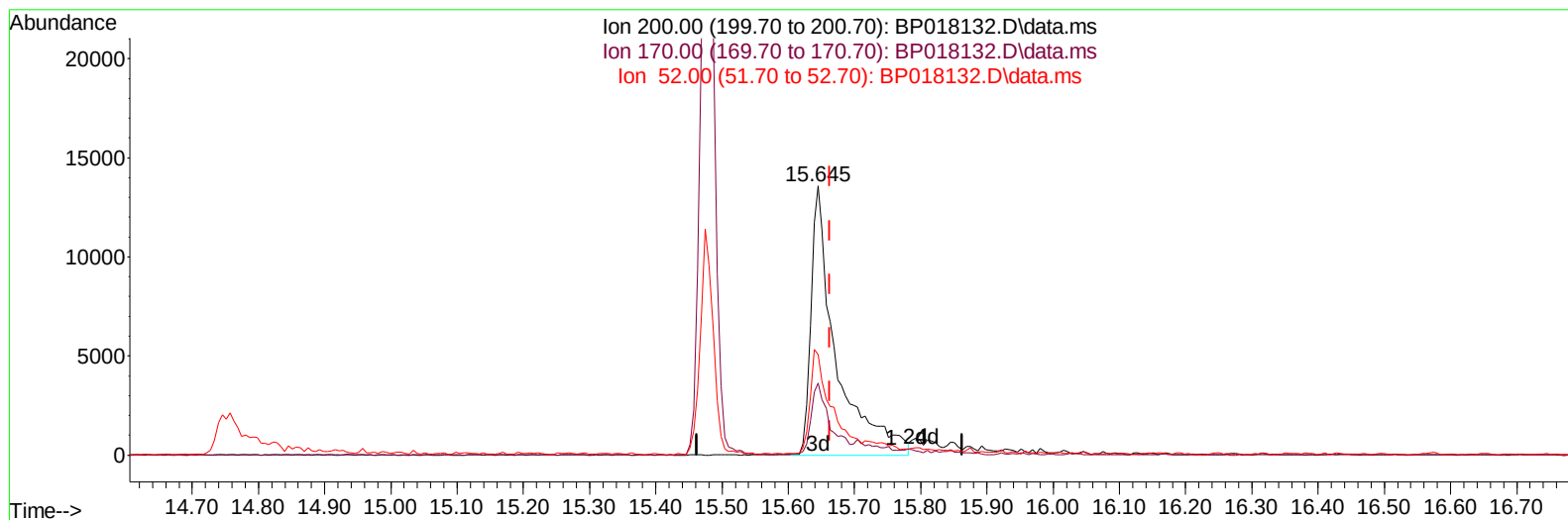
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TIC: BP018132.D\data.ms

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

15.645min (-0.018) 11.67 ng/ul m

response 35495

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	23.60	26.80
52.00	52.00	37.32#
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	7.787	152	69569	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	289730	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.475	164	186567	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.257	188	420731	20.000	ng/ul	0.00
79) Chrysene-d12	21.392	240	382909	20.000	ng/ul	0.00
88) Perylene-d12	23.868	264	405750	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.164	96	11297	5.728	ng/uL	0.00
4) Pyridine-d5	3.605	84	84952	16.521	ng/ul	0.00
7) Phenol-d5	6.958	99	166242	24.752	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.140	67	112434	28.486	ng/ul	0.00
11) 2-Chlorophenol-d4	7.311	132	136946	25.748	ng/ul	0.00
15) 4-Methylphenol-d8	8.516	113	136171	24.742	ng/ul	0.00
21) Nitrobenzene-d5	8.999	128	67526	25.869	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	76562	25.890	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.240	165	122631	23.267	ng/ul	0.00
31) 4-Chloroaniline-d4	10.793	131	142659	19.737	ng/ul	0.00
46) Dimethylphthalate-d6	13.898	166	489807	28.454	ng/ul	0.00
49) Acenaphthylene-d8	14.175	160	484073	25.818	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	40683	15.816	ng/ul	0.01
60) Fluorene-d10	15.475	176	385883	27.152	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	35495m	11.670	ng/ul	-0.02
73) Anthracene-d10	17.357	188	623234	27.389	ng/ul	0.00
81) Pyrene-d10	19.616	212	725020	28.347	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	672474	28.503	ng/ul	0.00

Target Compounds Qvalue

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

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