

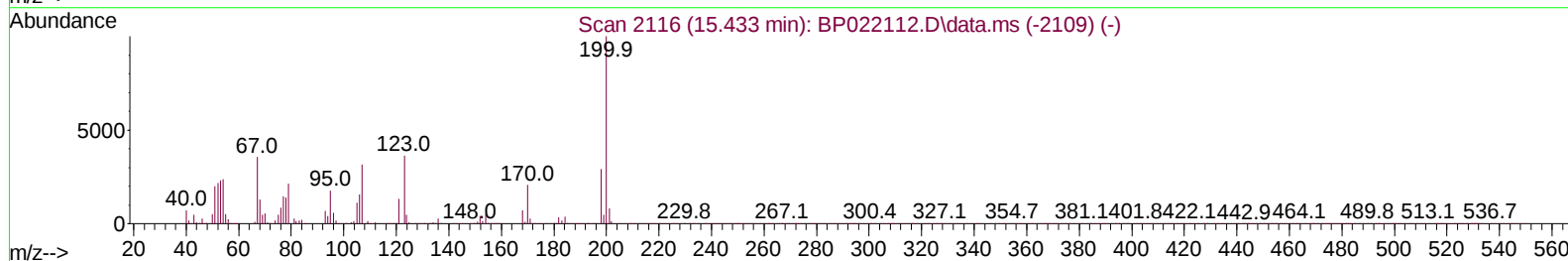
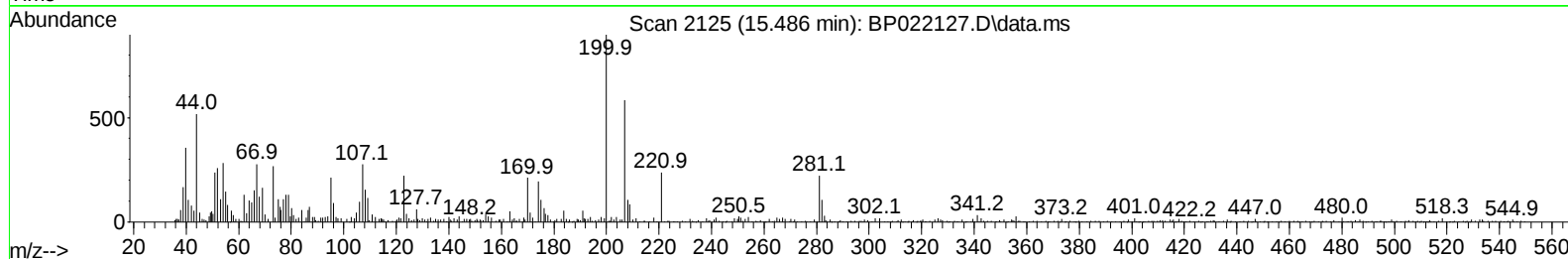
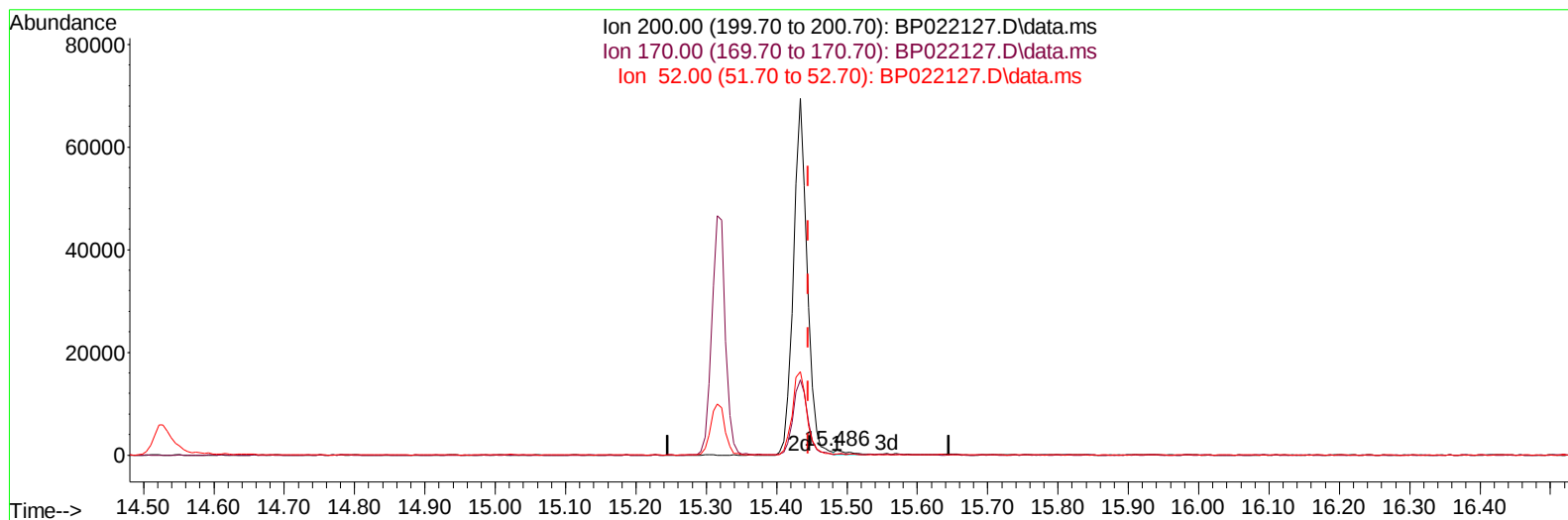
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP093024\
Data File : BP022127.D
Acq On : 01 Oct 2024 01:31
Operator : RC/JU
Sample : PB163565BL
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK565

Manual IntegrationsAPPROVED

Quant Time: Oct 01 02:16:44 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 24 15:23:13 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/01/2024
Supervised By :mohammad ahmed 10/02/2024



TIC: BP022127.D\data.ms

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

15.486min (+ 0.041) 0.23 ng/u1

response 1064

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.90	23.63
52.00	32.40	28.99
0.00	0.00	0.00

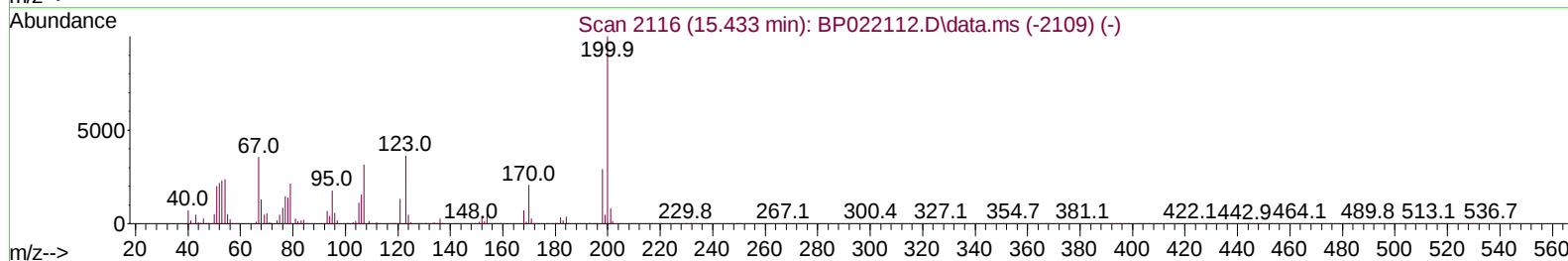
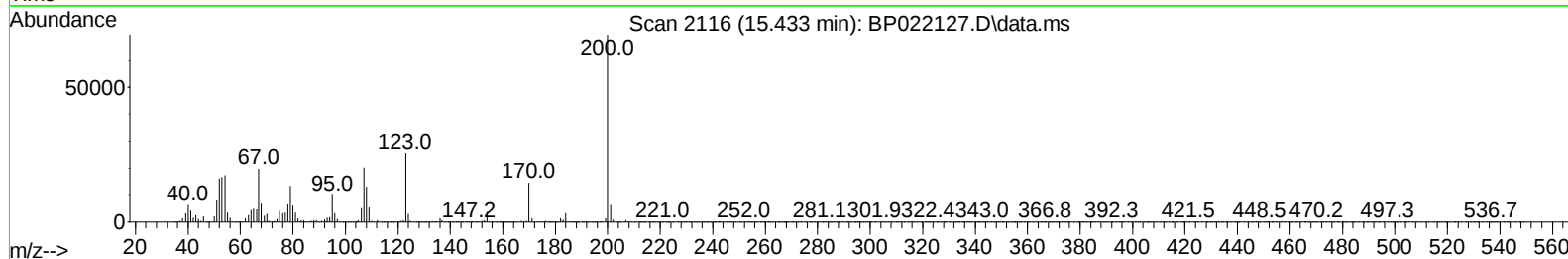
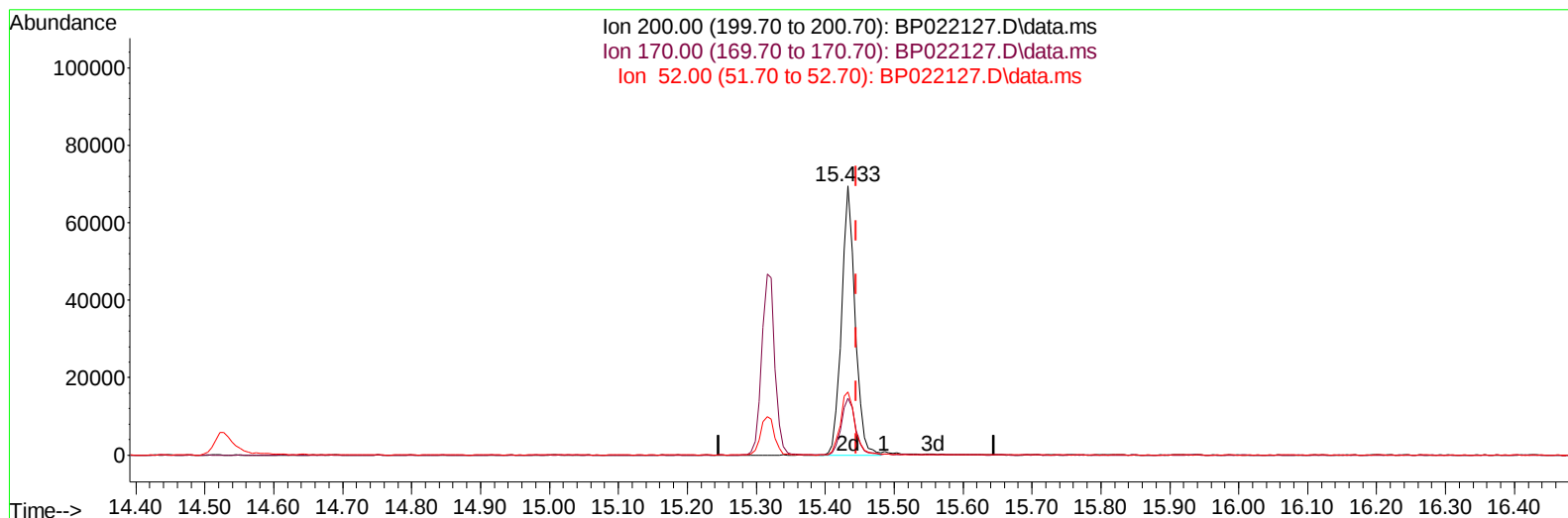
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TIC: BP022127.D\data.ms

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

15.433min (-0.012) 20.79 ng/ul m

response 94491

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	21.90	21.17
52.00	32.40	23.44#
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	7.699	152	112018	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.457	136	408580	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.310	164	307859	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.098	188	732055	20.000	ng/ul	-0.02
79) Chrysene-d12	21.533	240	854926	20.000	ng/ul	0.00
88) Perylene-d12	24.815	264	999063	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	14056	4.915	ng/uL	0.00
4) Pyridine-d5	3.634	84	186692	22.838	ng/ul	0.00
7) Phenol-d5	6.881	99	212208	23.159	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.040	67	134440	23.827	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	168178	25.367	ng/ul	0.00
15) 4-Methylphenol-d8	8.399	113	170667	23.505	ng/ul	0.00
21) Nitrobenzene-d5	8.840	128	81333	26.596	ng/ul	0.00
24) 2-Nitrophenol-d4	9.551	143	92276	26.613	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.087	165	181097	24.050	ng/ul	0.00
31) 4-Chloroaniline-d4	10.593	131	215544	23.991	ng/ul	-0.01
46) Dimethylphthalate-d6	13.710	166	606467	25.484	ng/ul	-0.01
49) Acenaphthylene-d8	13.998	160	624516	24.786	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143	85362	20.970	ng/ul	0.00
60) Fluorene-d10	15.316	176	514744	24.453	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.433	200	94491m	20.790	ng/ul	-0.01
73) Anthracene-d10	17.204	188	867960	25.397	ng/ul	-0.01
81) Pyrene-d10	19.574	212	1106479	24.981	ng/ul	-0.02
92) Benzo(a)pyrene-d12	24.592	264	1284568	24.446	ng/ul	-0.02

Target Compounds Qvalue

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

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