

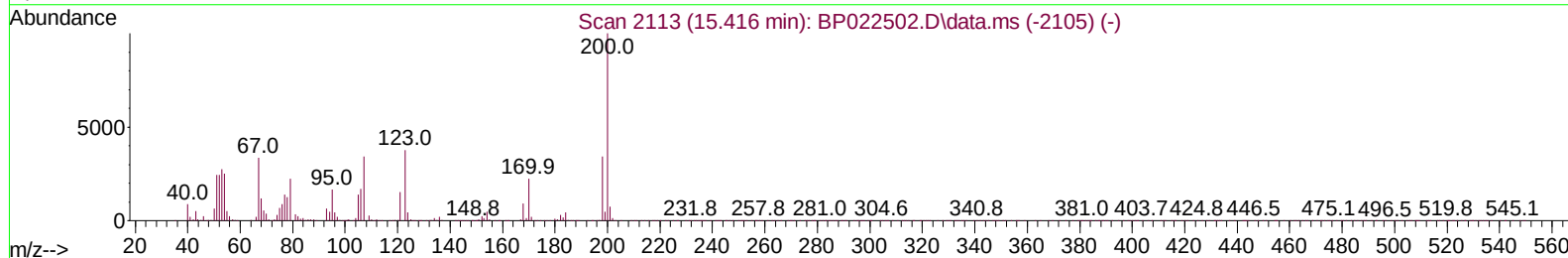
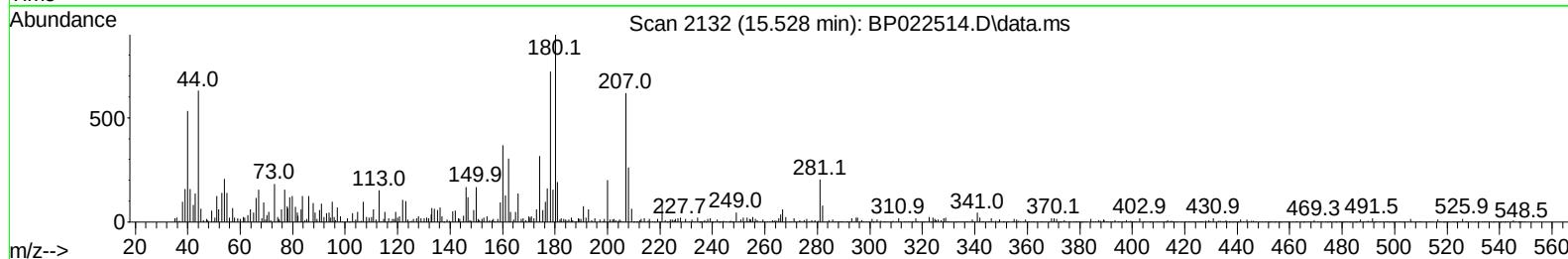
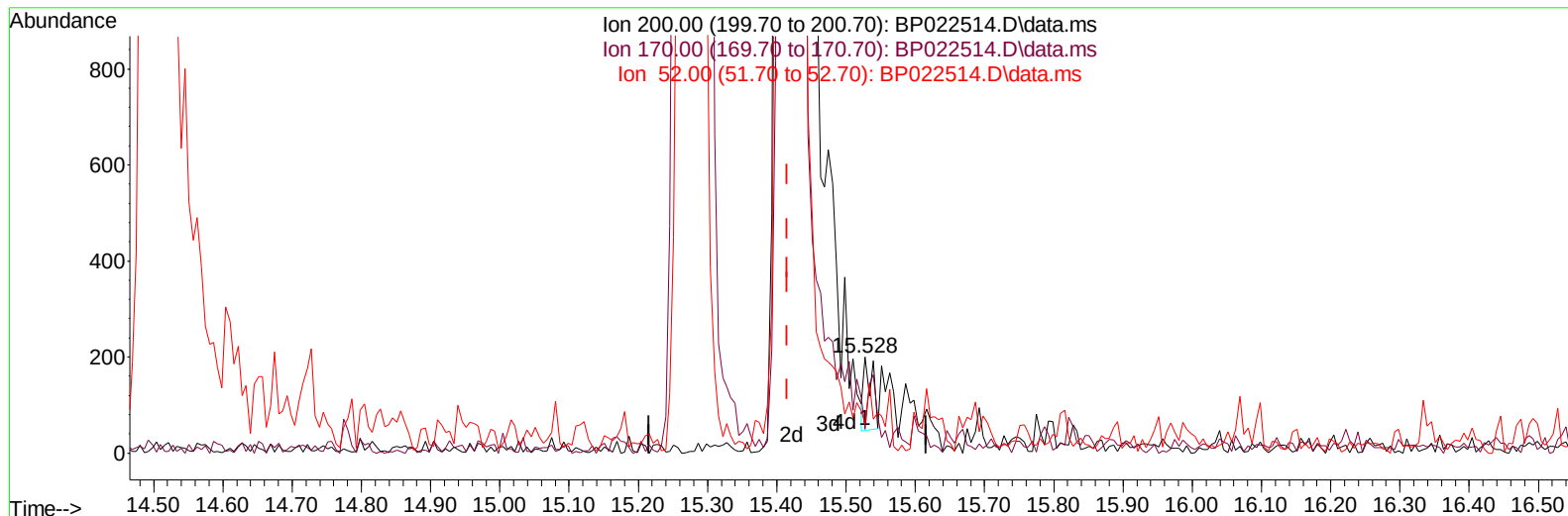
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102124\
Data File : BP022514.D
Acq On : 21 Oct 2024 20:52
Operator : RC/JU
Sample : PB164300BL
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK300

Manual IntegrationsAPPROVED

Quant Time: Oct 22 01:40:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 18 11:08:24 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/22/2024
Supervised By :mohammad ahmed 10/25/2024



TIC: BP022514.D\data.ms

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

15.528min (+ 0.112) 0.03 ng/u1

response 129

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.00	25.00
52.00	36.60	31.00
0.00	0.00	0.00

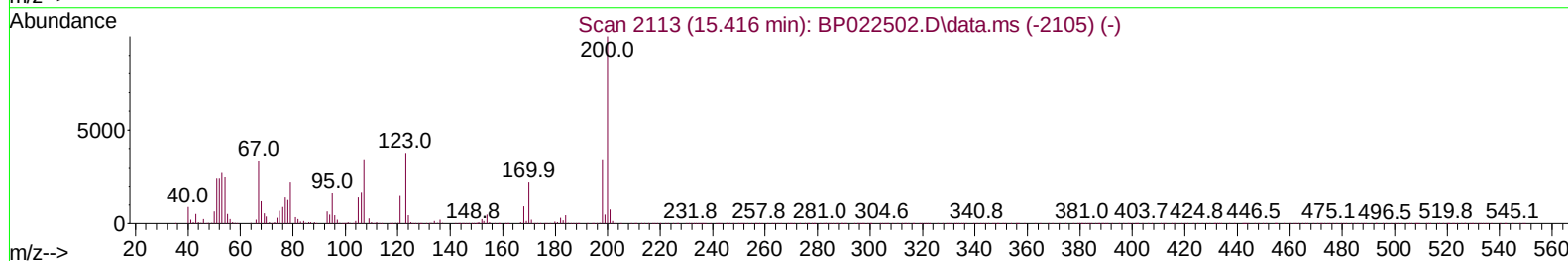
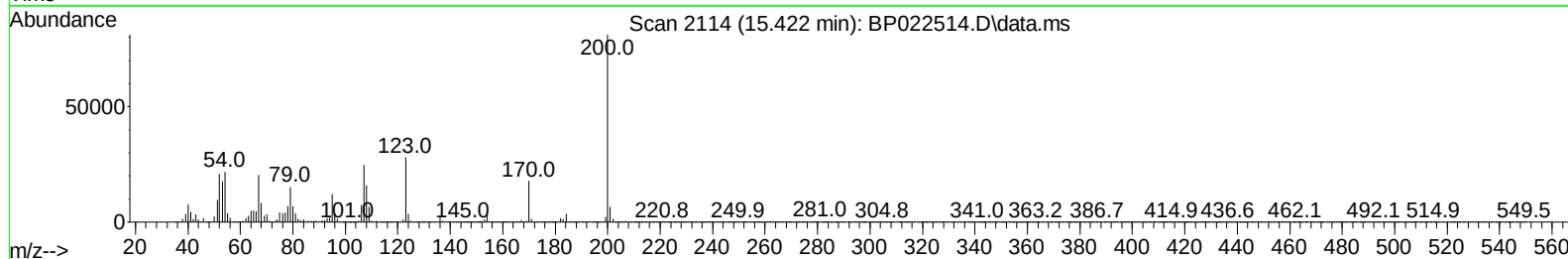
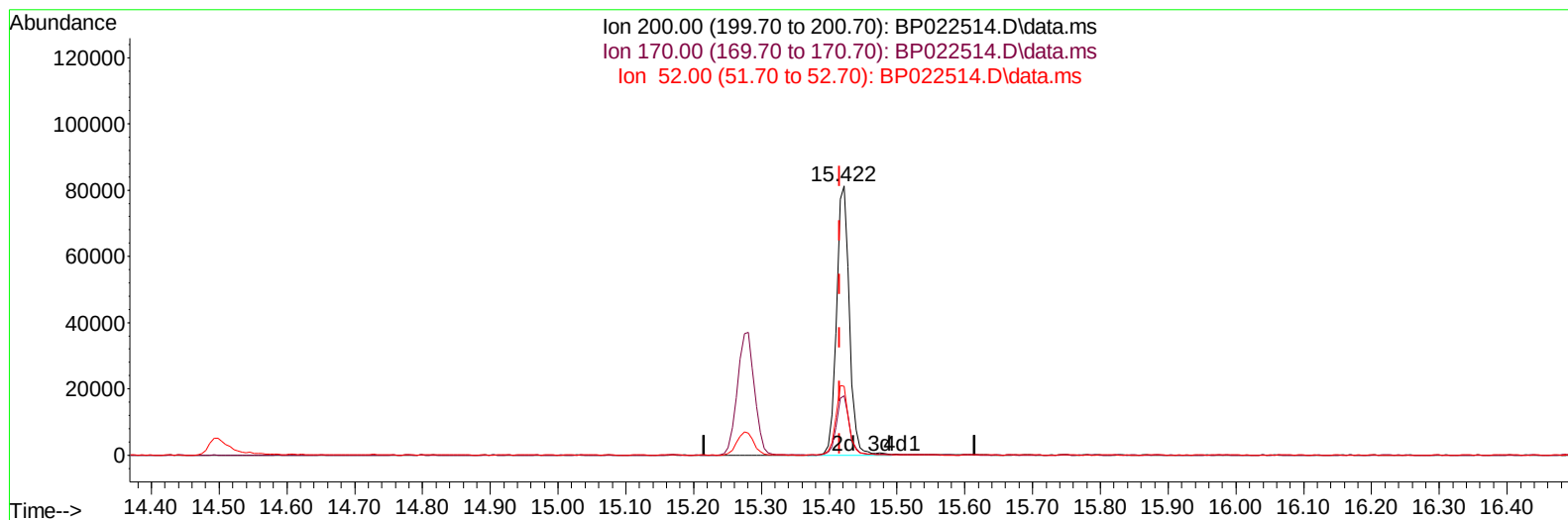
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.422min (+ 0.006) 22.86 ng/ul m

response 107363

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.00	21.96
52.00	36.60	25.64#
0.00	0.00	0.00

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Instrument :
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Reviewed By :Yogesh Patel 10/22/2024
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Quant Time: Oct 22 01:41:37 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.663	152	102173	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	381968	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	297351	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.080	188	713969	20.000	ng/ul	0.00
79) Chrysene-d12	21.504	240	803498	20.000	ng/ul	0.00
88) Perylene-d12	24.768	264	911432	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.199	96	12354	4.699	ng/uL	0.00
4) Pyridine-d5	3.611	84	163143	22.602	ng/ul	0.00
7) Phenol-d5	6.857	99	204502	23.778	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.005	67	118802	23.502	ng/ul	0.00
11) 2-Chlorophenol-d4	7.205	132	152870	25.049	ng/ul	0.00
15) 4-Methylphenol-d8	8.363	113	164656	23.989	ng/ul	0.00
21) Nitrobenzene-d5	8.804	128	77022	26.225	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	92957	27.338	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.051	165	180114	24.923	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	213616	25.015	ng/ul	0.00
46) Dimethylphthalate-d6	13.681	166	626152	26.504	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	644640	25.690	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	84275	21.263	ng/ul	0.00
60) Fluorene-d10	15.275	176	523529	25.548	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.422	200	107363m	22.864	ng/ul	0.00
73) Anthracene-d10	17.175	188	913945	27.211	ng/ul	0.00
81) Pyrene-d10	19.557	212	1170735	27.553	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.551	264	1308614	26.885	ng/ul	0.02

Target Compounds Qvalue

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

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