

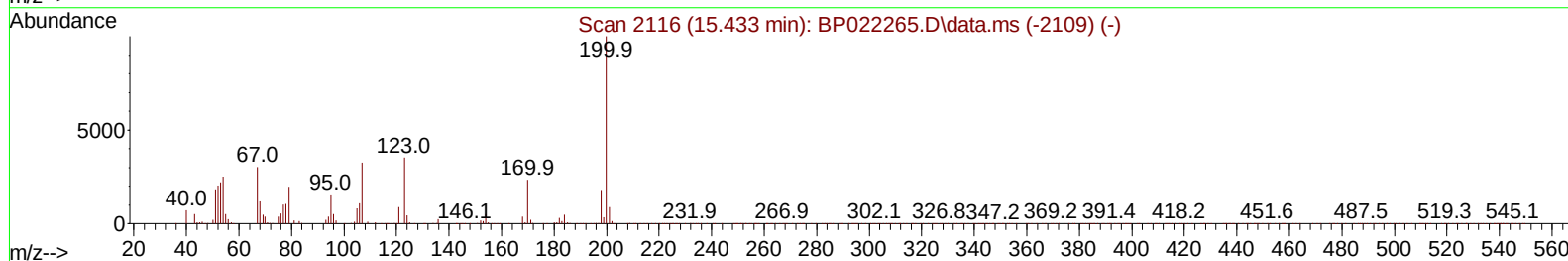
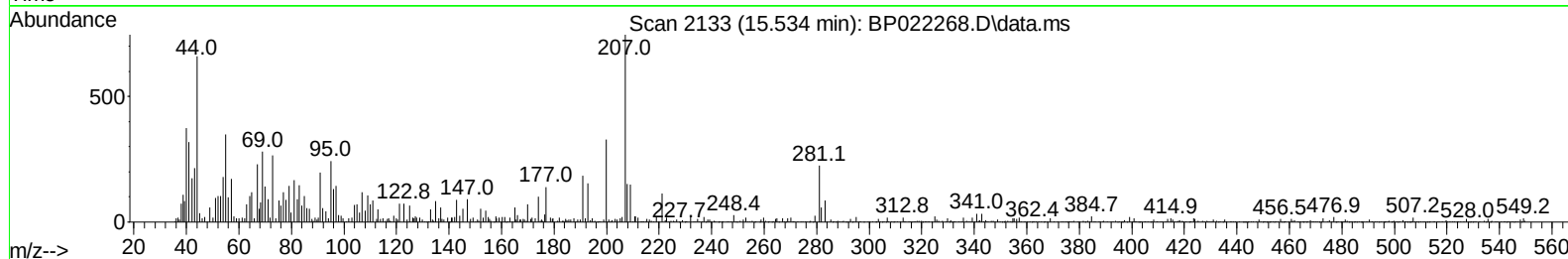
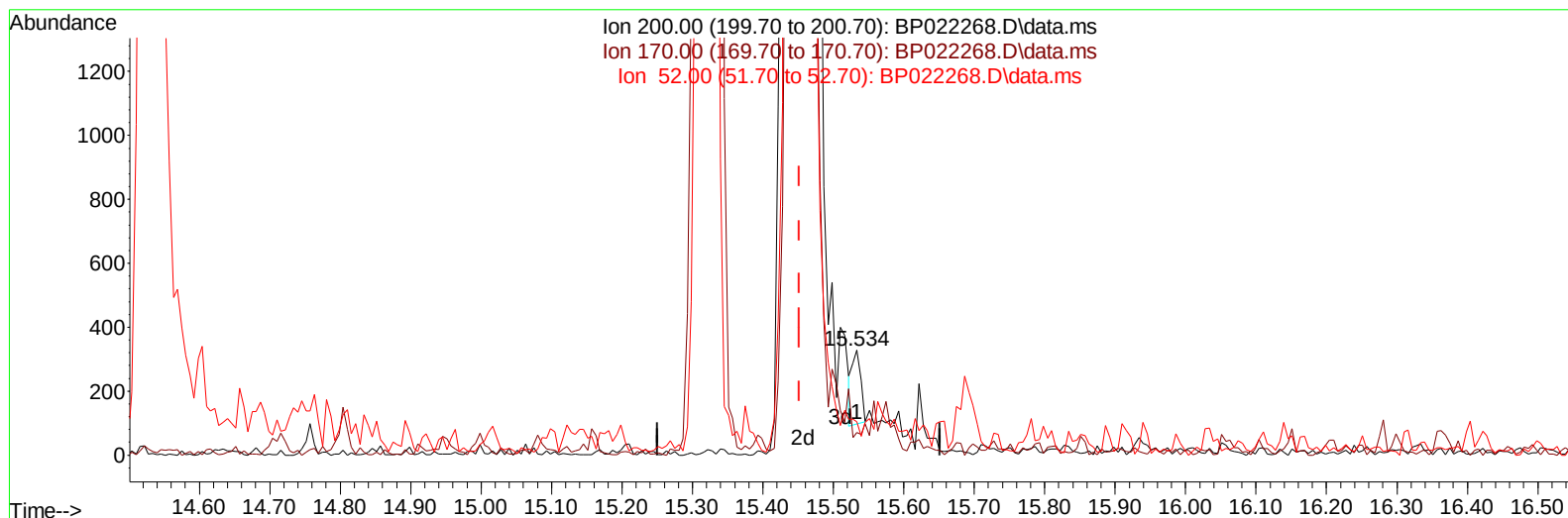
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100824\
Data File : BP022268.D
Acq On : 08 Oct 2024 12:42
Operator : RC/JU
Sample : P4162-04
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4ES0

Manual IntegrationsAPPROVED

Quant Time: Oct 08 13:08:09 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 08 02:14:47 2024
Response via : Initial Calibration

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



TIC: BP022268.D\data.ms

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

15.534min (+ 0.082) 0.03 ng/ul

response 199

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.00	21.34
52.00	36.60	32.01
0.00	0.00	0.00

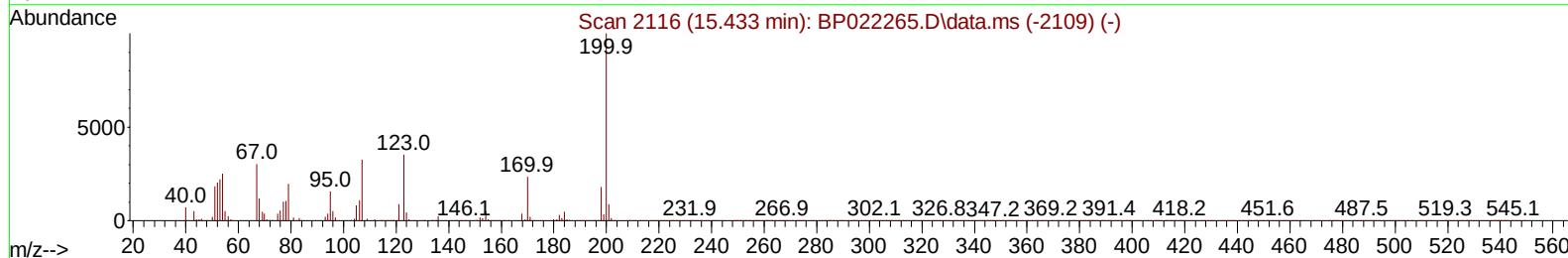
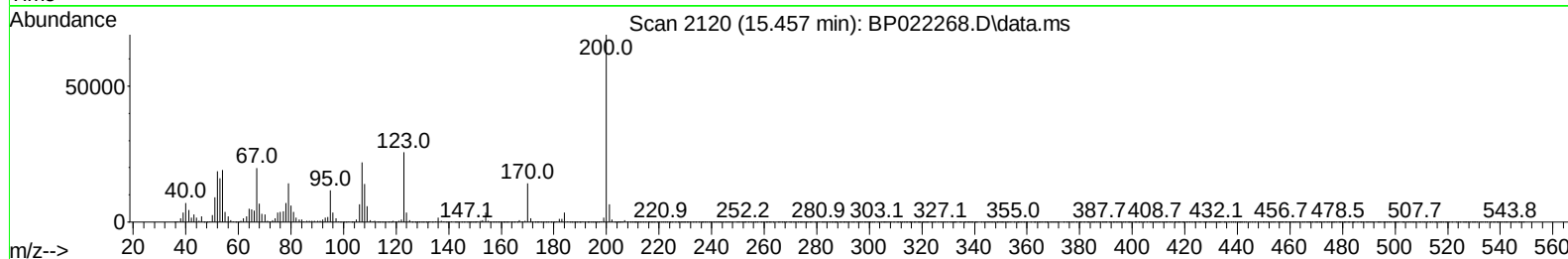
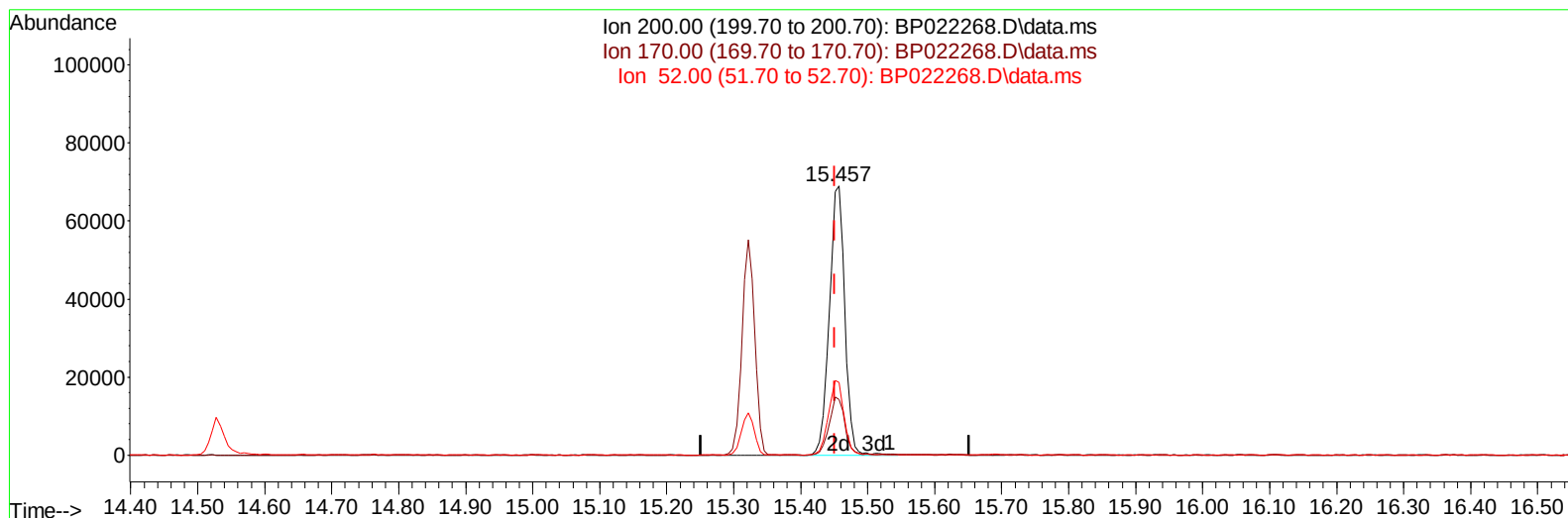
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.457min (+ 0.006) 15.82 ng/ul m

response 111016

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.00	20.71
52.00	36.60	27.04#
0.00	0.00	0.00

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Reviewed By :Yogesh Patel 10/10/2024
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Quant Time: Oct 08 13:10:41 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.705	152	148860	20.000	ng/ul	0.00
20) Naphthalene-d8	10.457	136	565889	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.310	164	452134	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.116	188	1067242	20.000	ng/ul	0.00
79) Chrysene-d12	21.539	240	1201450	20.000	ng/ul	0.00
88) Perylene-d12	24.827	264	1447253	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.228	96	10447	2.727	ng/uL	0.00
4) Pyridine-d5	3.640	84	151496	14.406	ng/ul	0.00
7) Phenol-d5	6.887	99	194220	15.500	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.040	67	121283	16.468	ng/ul	0.00
11) 2-Chlorophenol-d4	7.246	132	144236	16.222	ng/ul	0.00
15) 4-Methylphenol-d8	8.405	113	159552	15.955	ng/ul	0.00
21) Nitrobenzene-d5	8.846	128	72733	16.716	ng/ul	0.00
24) 2-Nitrophenol-d4	9.557	143	85476	16.968	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.087	165	172127	16.077	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.599	131	206341	16.310	ng/ul	0.00
46) Dimethylphthalate-d6	13.722	166	698476	19.444	ng/ul	0.00
49) Acenaphthylene-d8	14.004	160	684693	17.945	ng/ul	0.00
54) 4-Nitrophenol-d4	14.528	143	97972	16.257	ng/ul	0.00
60) Fluorene-d10	15.322	176	592836	19.026	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.457	200	111016m	15.816	ng/ul	0.00
73) Anthracene-d10	17.210	188	987927	19.678	ng/ul	0.00
81) Pyrene-d10	19.586	212	1276292	20.088	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.610	264	1573699	20.361	ng/ul	0.01

Target Compounds Qvalue

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

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