

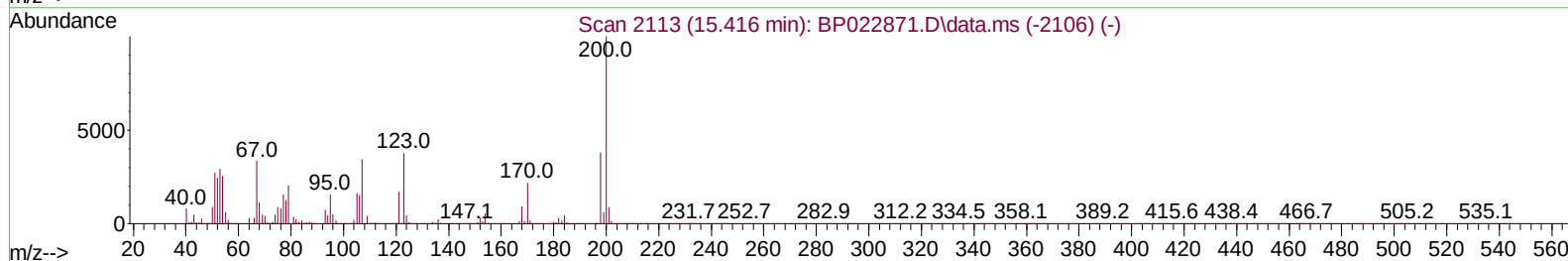
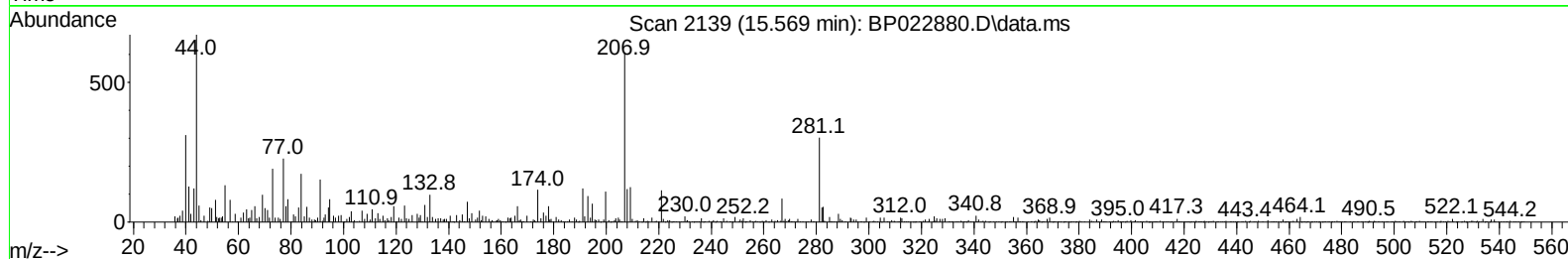
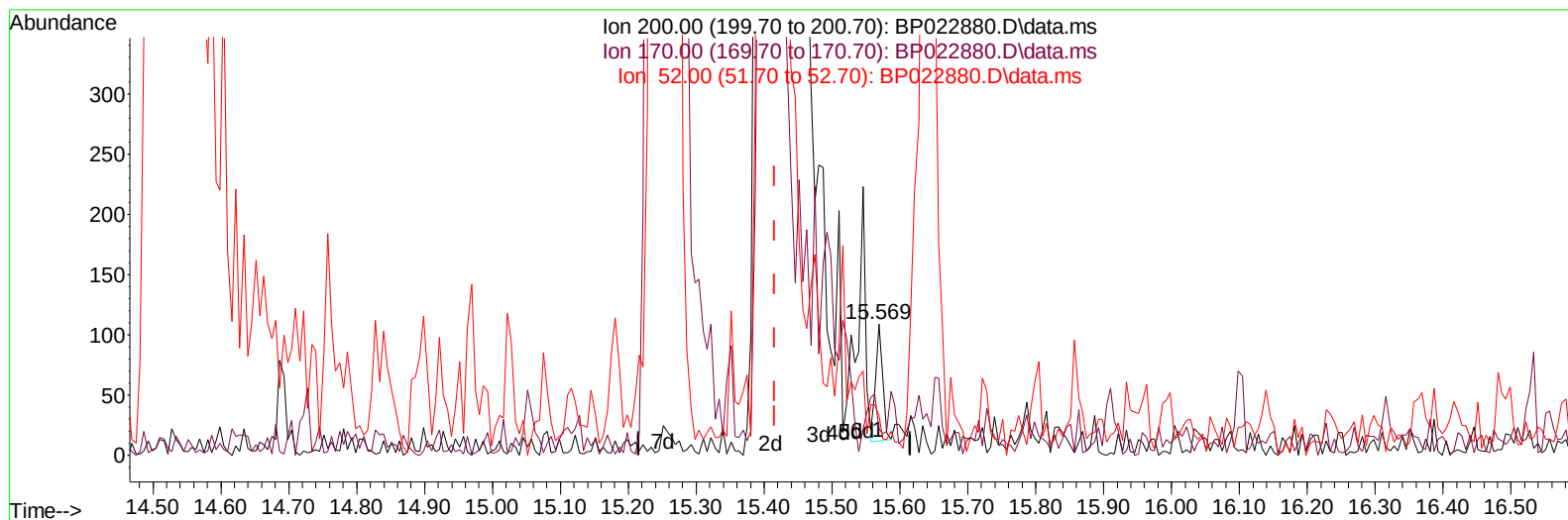
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022880.D
Acq On : 06 Nov 2024 05:01
Operator : RC/JU
Sample : P4581-16
Misc :
ALS Vial : 26 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A4ED7

Manual IntegrationsAPPROVED

Quant Time: Nov 06 05:26:01 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 06 00:31:03 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/06/2024
Supervised By :mohammad ahmed 11/08/2024



TIC: BP022880.D\data.ms

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

15.569min (+ 0.153) 0.01 ng/u1

response 69

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.00	20.18
52.00	36.60	30.28
0.00	0.00	0.00

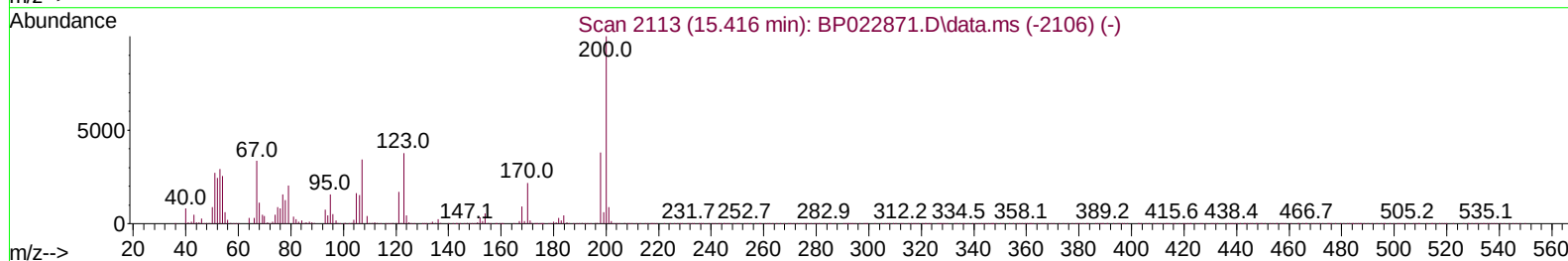
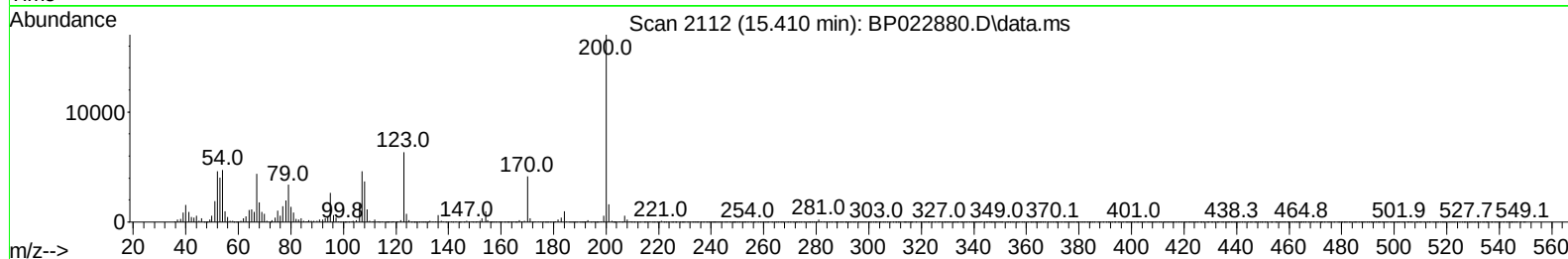
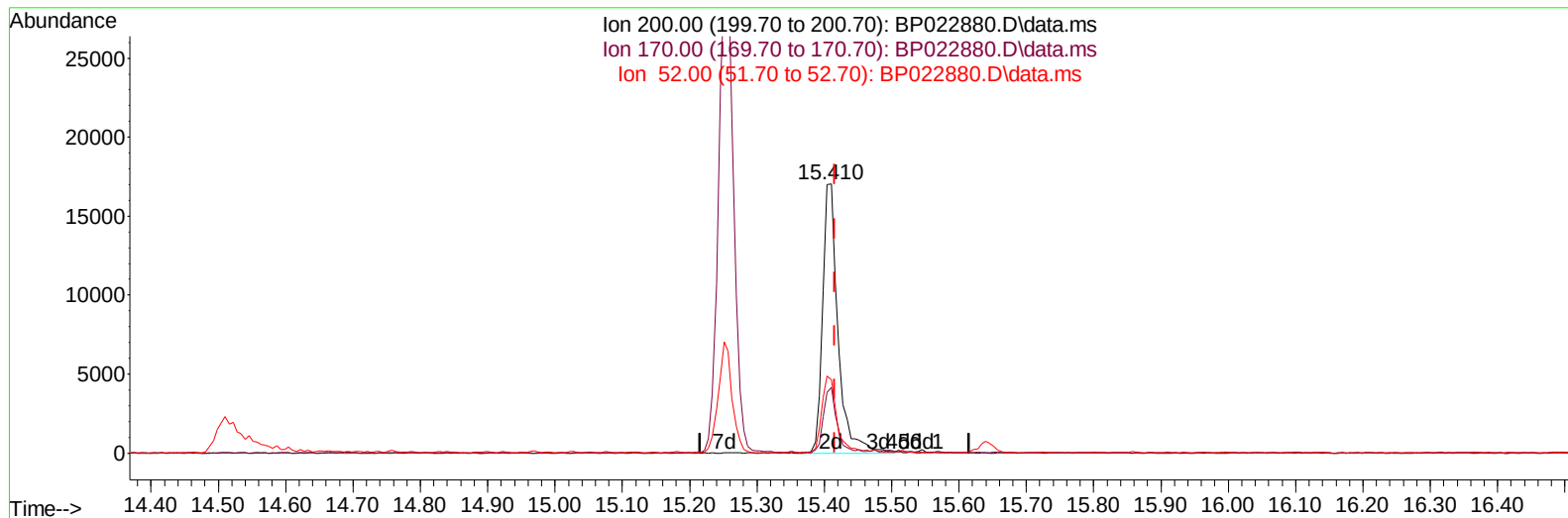
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(65) 4,6-Dinitro-2-methylphenol-d2 (S)

15.410min (-0.006) 5.60 ng/ul m

response 26914

Ion	Exp%	Act%
200.00	100.00	100.00
170.00	22.00	24.36
52.00	36.60	27.21#
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	7.622	152	101425	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	386616	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.245	164	298350	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.051	188	730955	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	769838	20.000	ng/ul	0.00
88) Perylene-d12	24.739	264	895949	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	5455	2.090	ng/uL	0.00
4) Pyridine-d5	3.593	84	64641	9.022	ng/ul	0.00
7) Phenol-d5	6.822	99	144037	16.871	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.969	67	79433	15.830	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	104859	17.309	ng/ul	0.00
15) 4-Methylphenol-d8	8.334	113	120658	17.709	ng/ul	0.00
21) Nitrobenzene-d5	8.763	128	52195	17.558	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	65146	18.929	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.016	165	138139	18.885	ng/ul	0.00
31) 4-Chloroaniline-d4	10.522	131	151028	17.474	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166	416702	17.579	ng/ul	0.00
49) Acenaphthylene-d8	13.934	160	485370	19.278	ng/ul	0.00
54) 4-Nitrophenol-d4	14.510	143	50589	12.721	ng/ul	0.01
60) Fluorene-d10	15.257	176	391919	19.062	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	26914m	5.598	ng/ul	0.00
73) Anthracene-d10	17.151	188	636102	18.499	ng/ul	-0.01
81) Pyrene-d10	19.539	212	796742	19.571	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.527	264	896375	18.734	ng/ul	0.00

Target Compounds Qvalue

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

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