

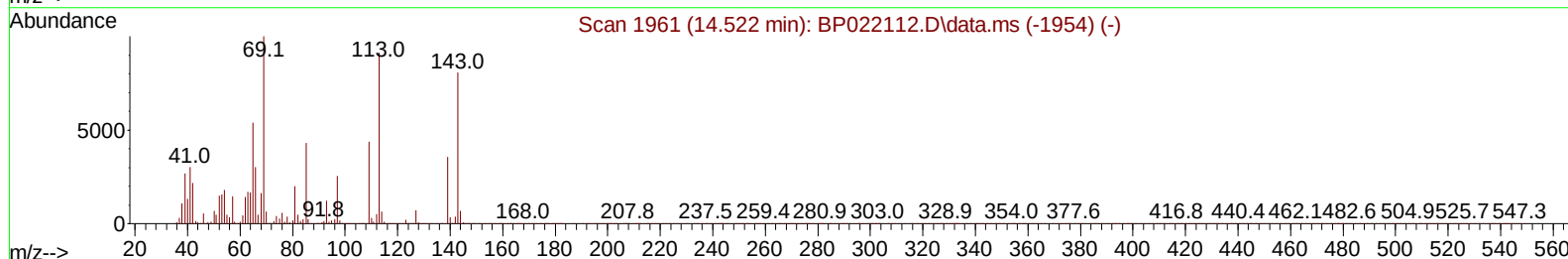
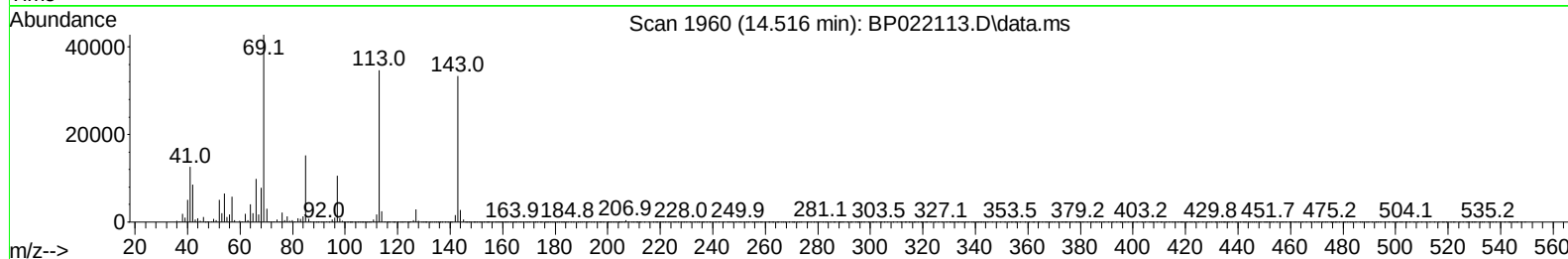
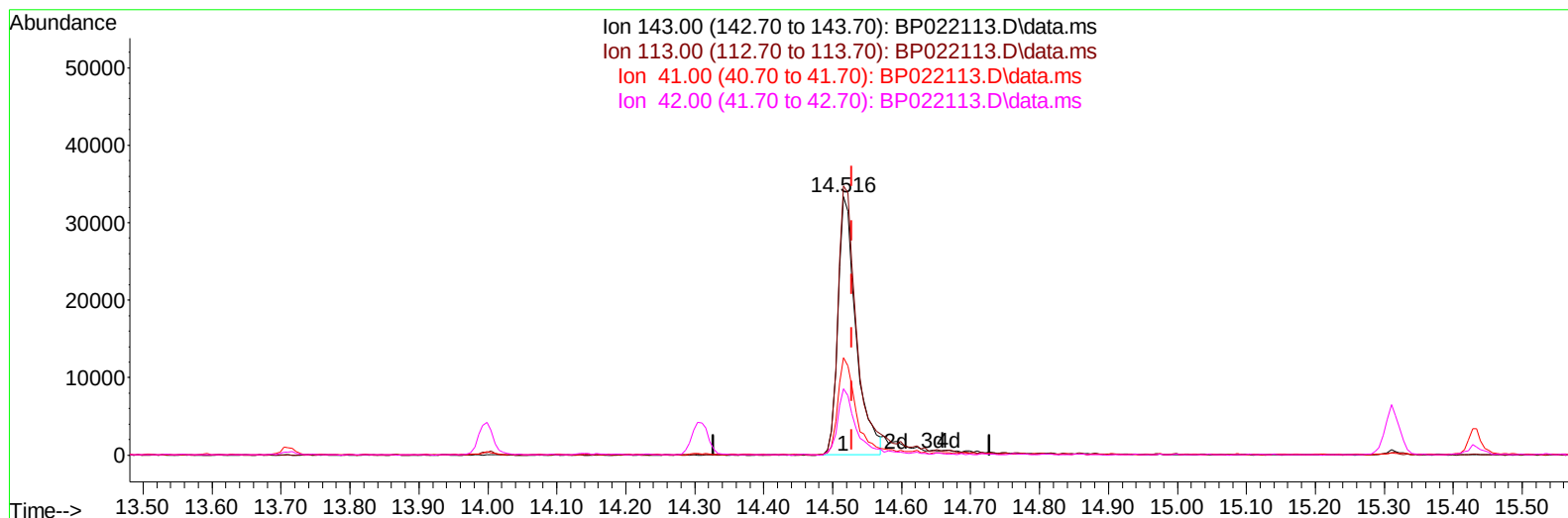
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP093024\
 Data File : BP022113.D
 Acq On : 30 Sep 2024 15:19
 Operator : RC/JU
 Sample : PB163717BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK717

Manual IntegrationsAPPROVED

Quant Time: Sep 30 16:30:53 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: BP022113.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (-0.012) 19.49 ng/ul

response 60538

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	104.20
41.00	35.80	37.82
42.00	26.20	25.61

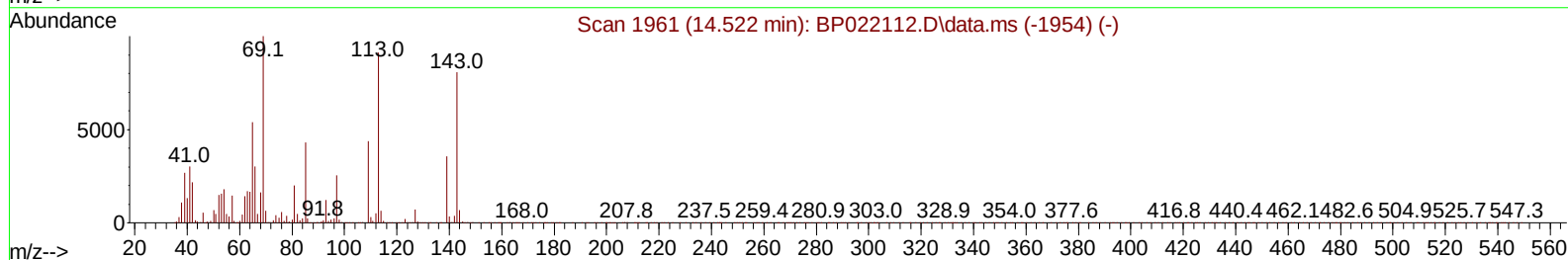
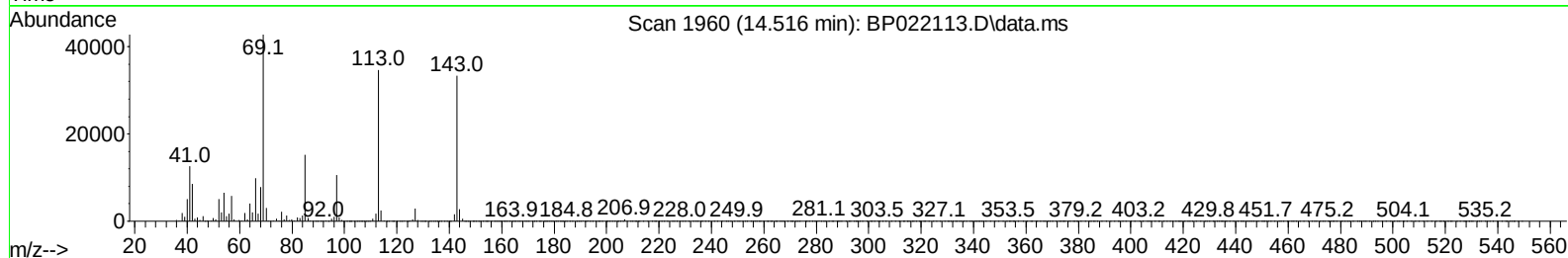
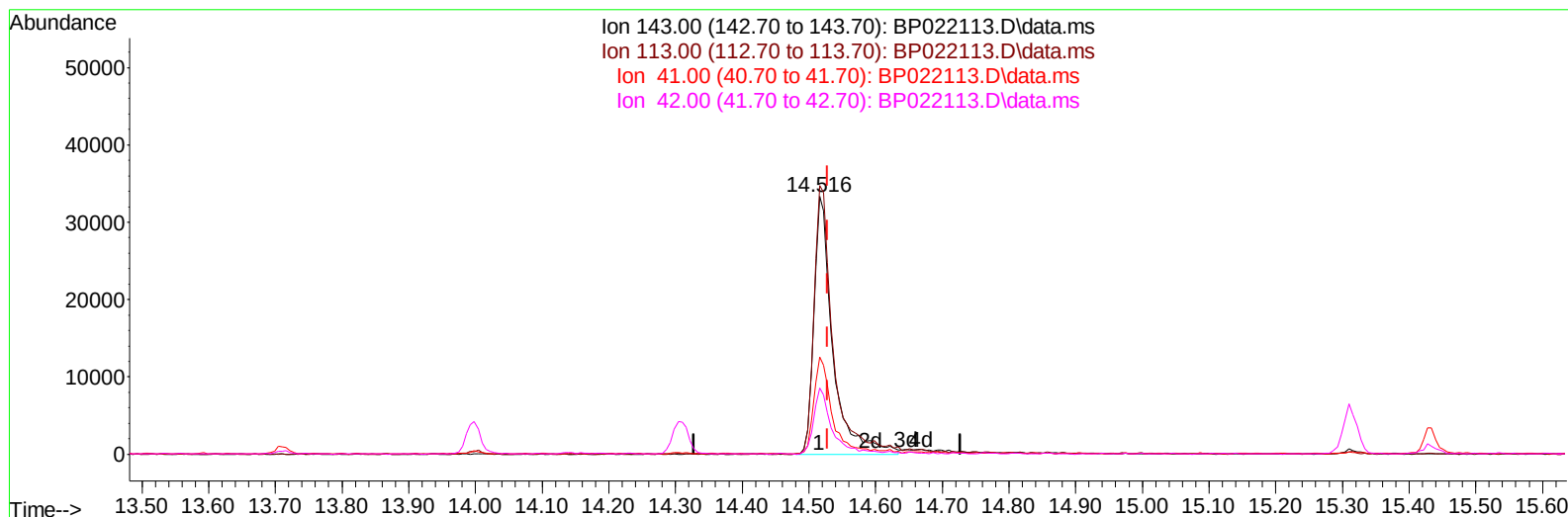
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TIC: BP022113.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.516min (-0.012) 21.12 ng/ul m

response 65619

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	106.70	104.20
41.00	35.80	37.82
42.00	26.20	25.61

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.699	152	93009	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.457	136	327850	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.310	164	234960	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.104	188	534392	20.000	ng/ul	-0.01
79) Chrysene-d12	21.533	240	605126	20.000	ng/ul	0.00
88) Perylene-d12	24.815	264	728380	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.222	96	13846	5.831	ng/uL	0.00
4) Pyridine-d5	3.634	84	173457	25.556	ng/ul	0.00
7) Phenol-d5	6.881	99	188068	24.719	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.040	67	118634	25.323	ng/ul	0.00
11) 2-Chlorophenol-d4	7.240	132	149536	27.165	ng/ul	0.00
15) 4-Methylphenol-d8	8.393	113	147922	24.536	ng/ul	-0.01
21) Nitrobenzene-d5	8.840	128	68251	27.814	ng/ul	0.00
24) 2-Nitrophenol-d4	9.552	143	80326	28.871	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.087	165	155084	25.667	ng/ul	0.00
31) 4-Chloroaniline-d4	10.593	131	185976	25.797	ng/ul	-0.01
46) Dimethylphthalate-d6	13.710	166	504844	27.796	ng/ul	-0.01
49) Acenaphthylene-d8	13.998	160	531254	27.627	ng/ul	0.00
54) 4-Nitrophenol-d4	14.516	143	65619m	21.122	ng/ul	-0.01
60) Fluorene-d10	15.310	176	423803	26.379	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.434	200	74063	22.323	ng/ul	-0.01
73) Anthracene-d10	17.198	188	712349	28.554	ng/ul	-0.02
81) Pyrene-d10	19.586	212	895343	28.559	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.592	264	1072190	27.987	ng/ul	-0.02

Target Compounds Qvalue

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

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