

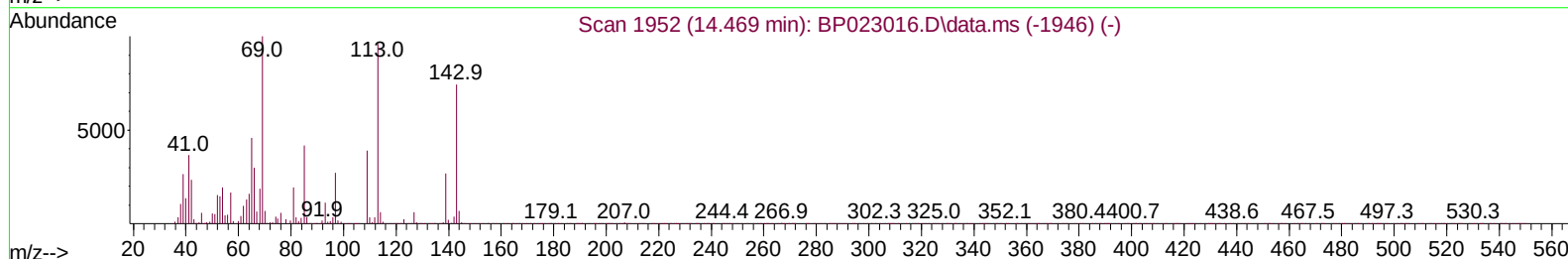
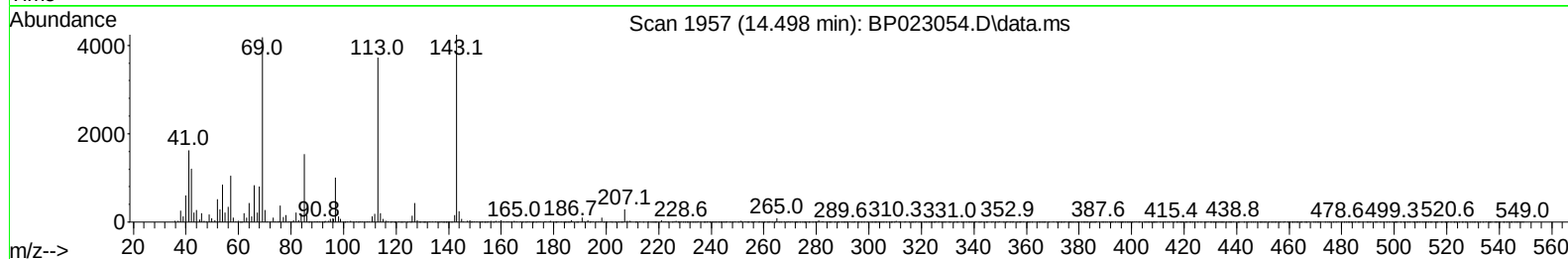
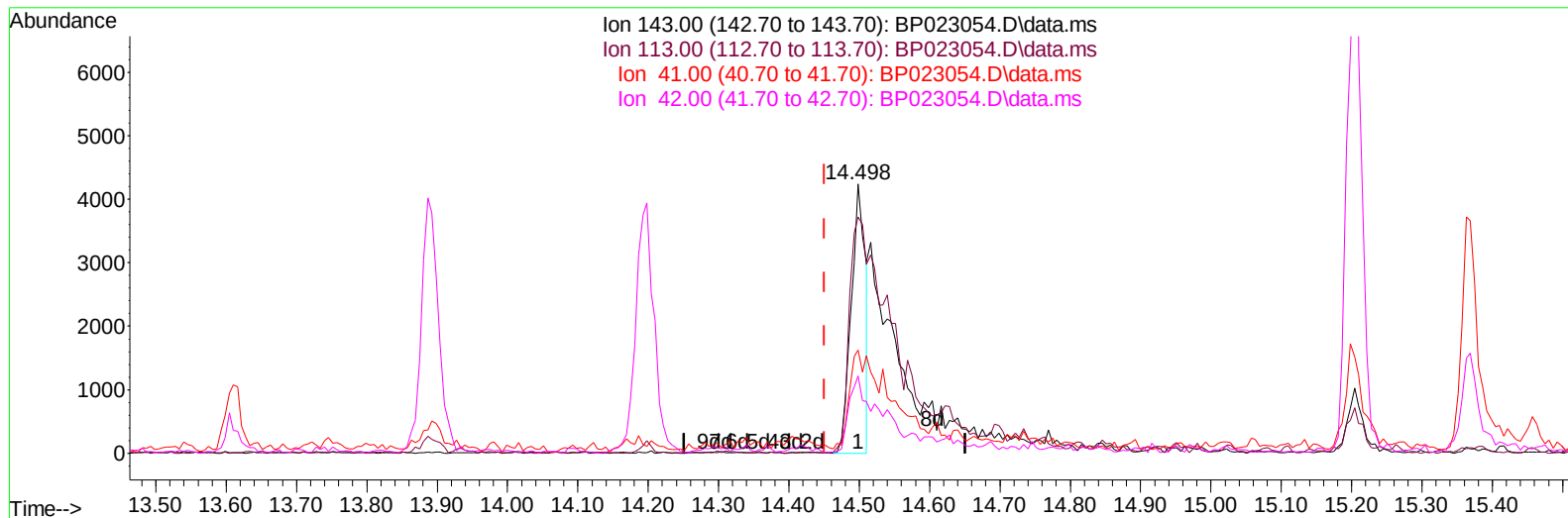
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111224\
Data File : BP023054.D
Acq On : 12 Nov 2024 19:58
Operator : RC/JU
Sample : P4687-03
Misc :
ALS Vial : 18 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0CN8

Manual IntegrationsAPPROVED

Quant Time: Nov 13 00:02:19 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Nov 12 23:59:19 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/14/2024
Supervised By :mohammad ahmed 11/14/2024



TIC: BP023054.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.047) 2.41 ng/ul

response 5941

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	87.85
41.00	39.20	38.42
42.00	25.90	28.56

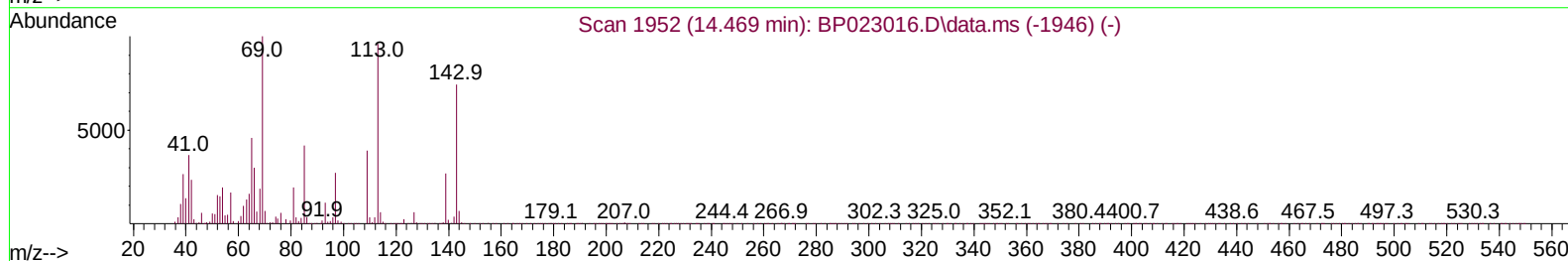
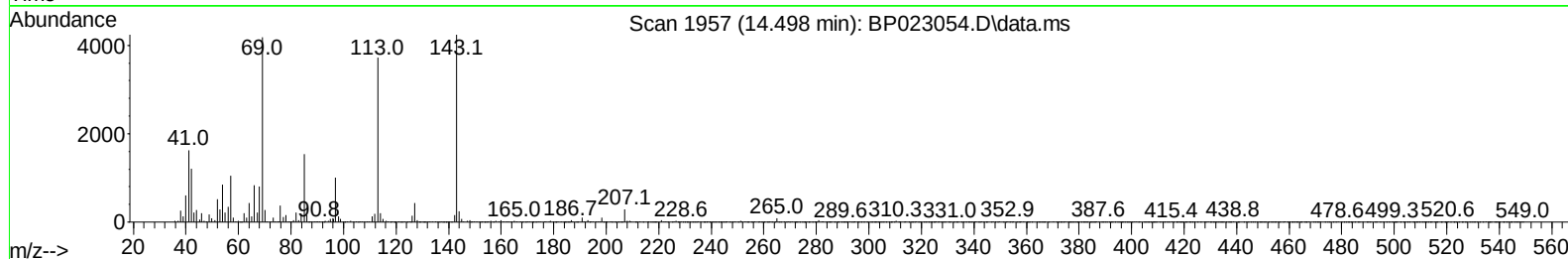
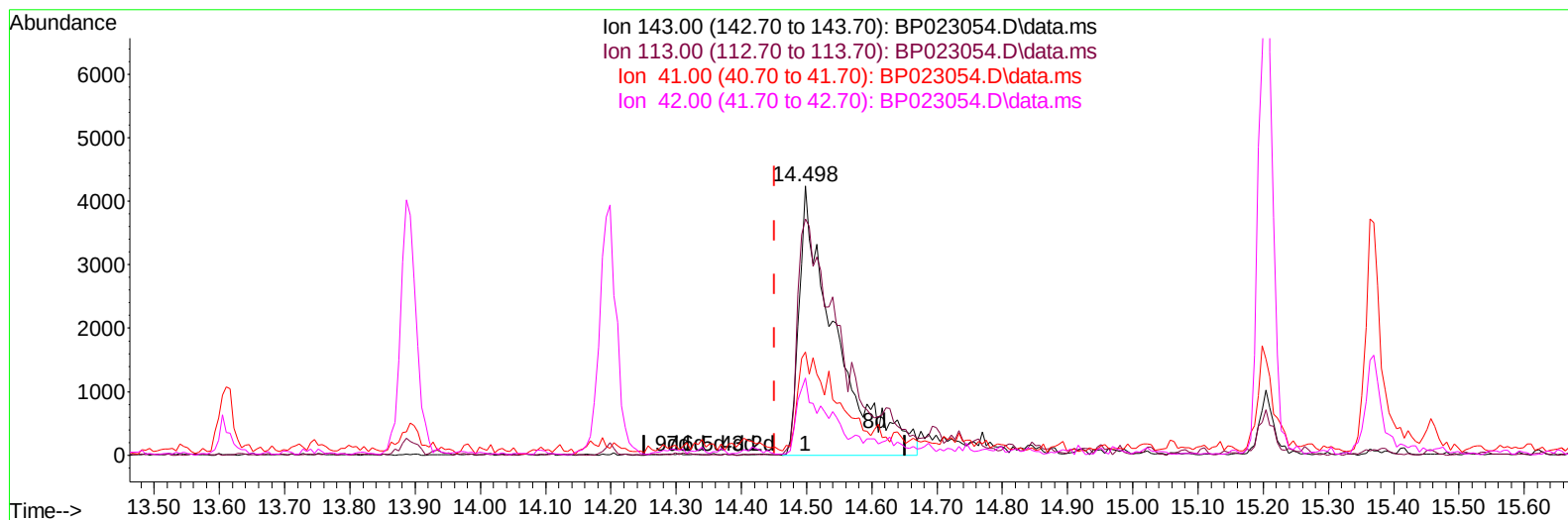
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TIC: BP023054.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.047) 6.64 ng/ul m

response 16354

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	87.85
41.00	39.20	38.42
42.00	25.90	28.56

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/14/2024
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Quant Time: Nov 13 02:14:28 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 12 23:59:19 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.581	152	71340	20.000	ng/ul	0.00
20) Naphthalene-d8	10.328	136	278985	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.198	164	227229	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.998	188	584469	20.000	ng/ul	-0.01
79) Chrysene-d12	21.451	240	659121	20.000	ng/ul	0.00
88) Perylene-d12	24.662	264	787743	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	7831	5.276	ng/uL	0.00
4) Pyridine-d5	3.558	84	41998	9.524	ng/ul	0.00
7) Phenol-d5	6.799	99	41175	7.807	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.928	67	96349	31.092	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	104041	26.934	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113	84743	19.525	ng/ul	0.00
21) Nitrobenzene-d5	8.728	128	59026	30.391	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143	71239	30.767	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	144945	29.892	ng/ul	0.00
31) 4-Chloroaniline-d4	10.481	131	155984	26.964	ng/ul	0.00
46) Dimethylphthalate-d6	13.610	166	541152	33.043	ng/ul	0.00
49) Acenaphthylene-d8	13.892	160	515377	30.160	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	16354m	6.645	ng/ul	0.05
60) Fluorene-d10	15.204	176	461490	32.140	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.369	200	100018	28.566	ng/ul	0.00
73) Anthracene-d10	17.110	188	868816	35.108	ng/ul	-0.01
81) Pyrene-d10	19.492	212	1142508	37.633	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.433	264	1371595	36.553	ng/ul	-0.03

Target Compounds Qvalue

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

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