

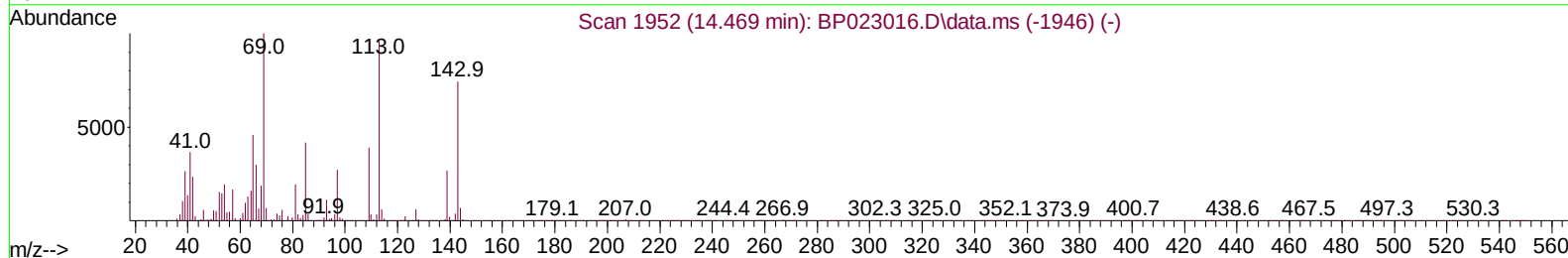
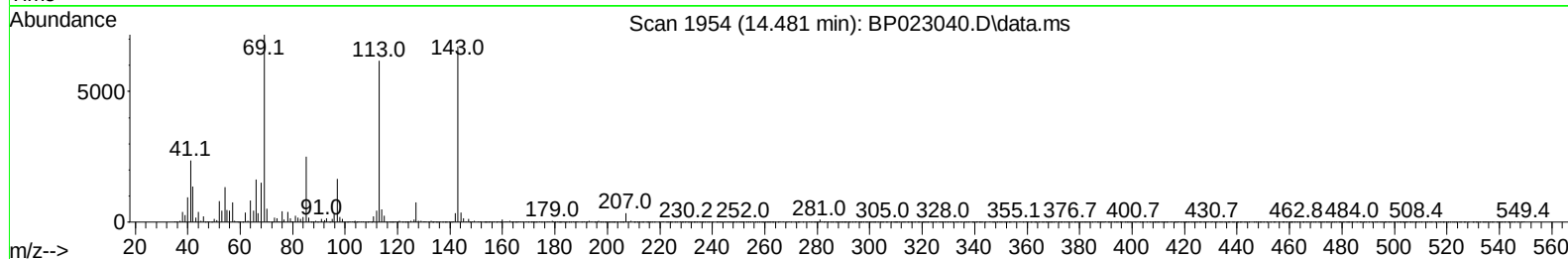
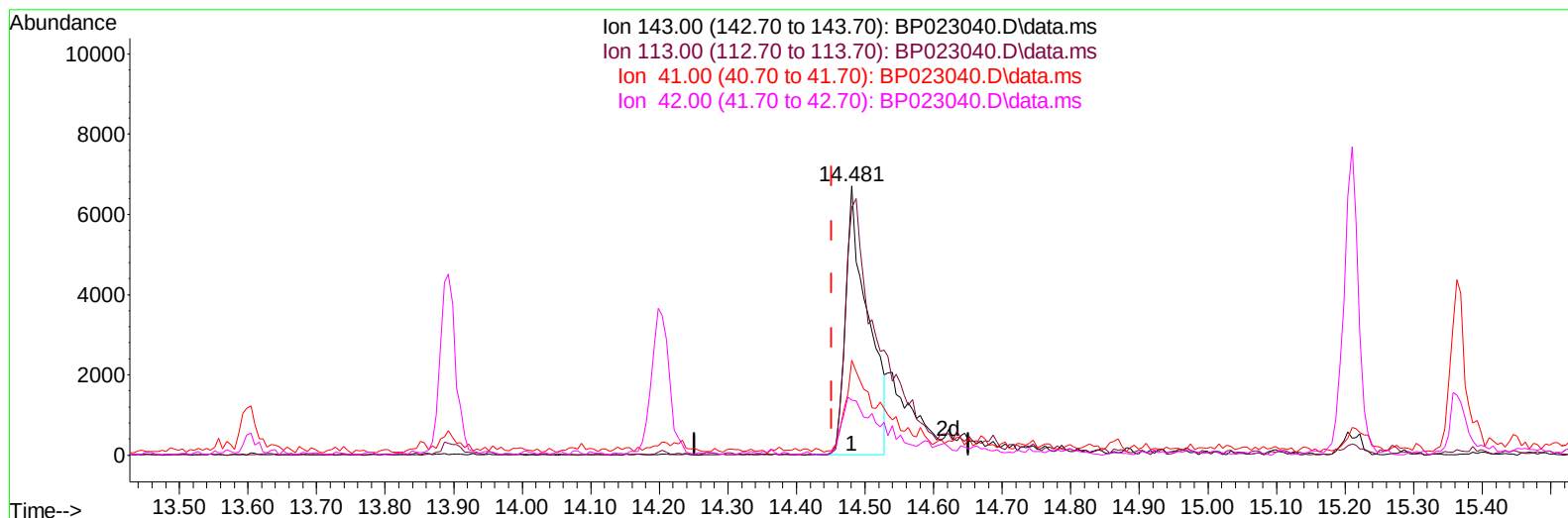
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111224\
Data File : BP023040.D
Acq On : 12 Nov 2024 10:30
Operator : RC/JU
Sample : P4761-18
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A0AN4

Manual IntegrationsAPPROVED

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

Quant Time: Nov 13 00:33:39 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



TIC: BP023040.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.029) 6.57 ng/u1

response 14924

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	92.02
41.00	39.20	35.34
42.00	25.90	20.37#

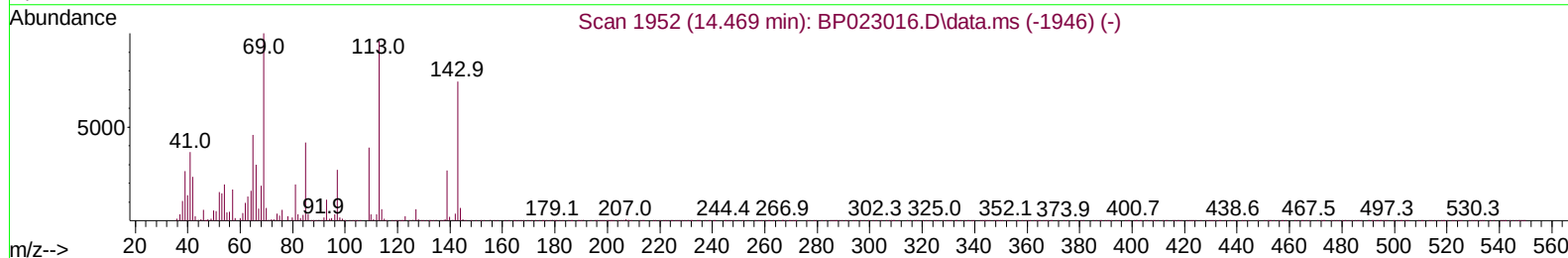
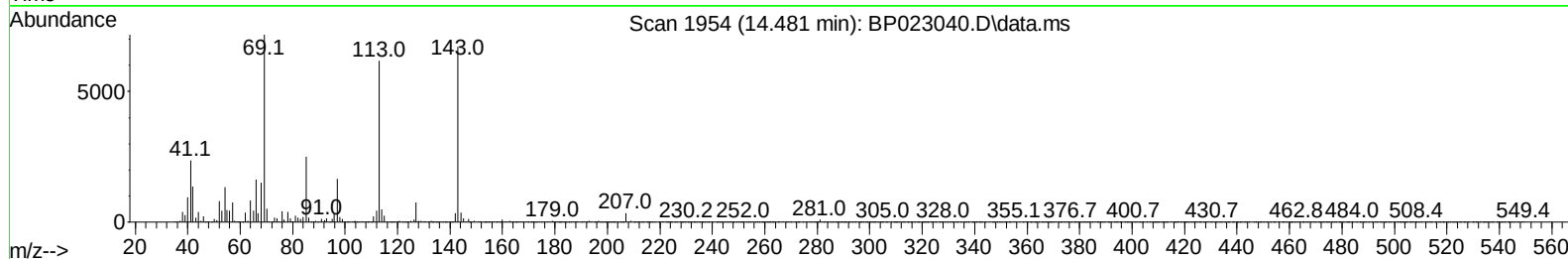
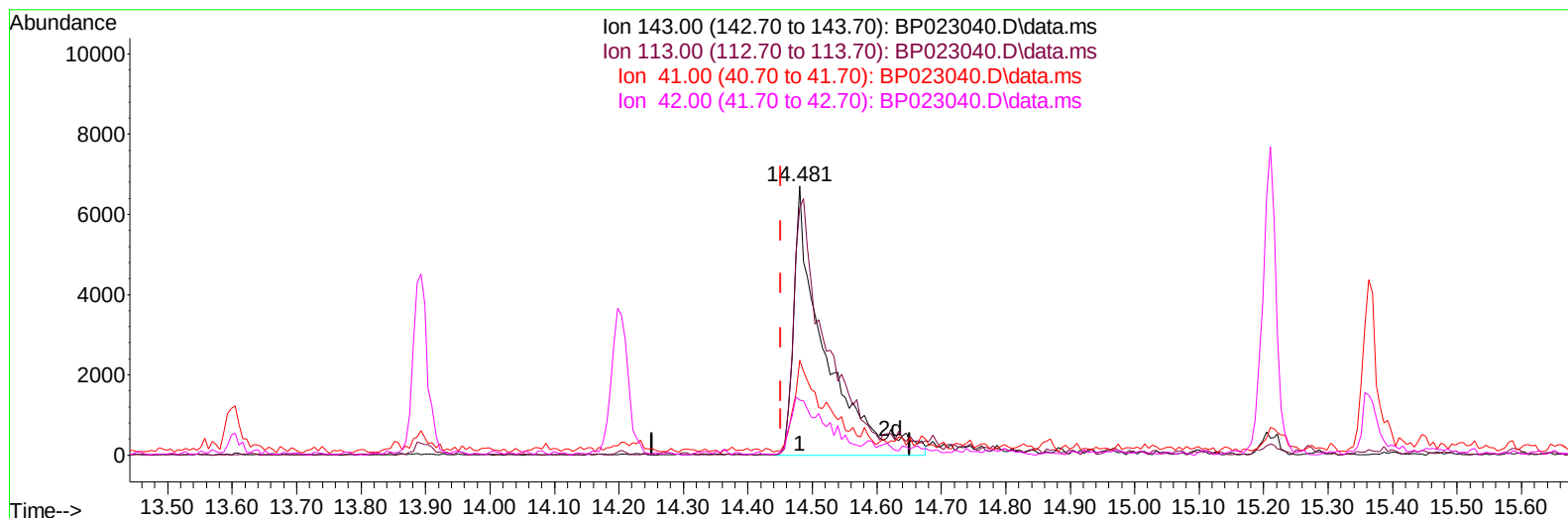
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TIC: BP023040.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.481min (+ 0.029) 9.64 ng/ul m

response 21910

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	92.02
41.00	39.20	35.34
42.00	25.90	20.37#

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Reviewed By :Yogesh Patel 11/14/2024
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Quant Time: Nov 13 00:46:15 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.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.587	152		66914	20.000	ng/ul	0.00
20) Naphthalene-d8	10.334	136		256114	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.204	164		209732	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.004	188		519911	20.000	ng/ul	0.00
79) Chrysene-d12	21.445	240		616599	20.000	ng/ul	0.00
88) Perylene-d12	24.674	264		750029	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.140	96		7368	5.293	ng/uL	0.00
4) Pyridine-d5	3.558	84		58111	14.050	ng/ul	0.00
7) Phenol-d5	6.799	99		57264	11.575	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.934	67		96590	33.232	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132		114162	31.509	ng/ul	0.00
15) 4-Methylphenol-d8	8.293	113		103038	25.310	ng/ul	0.00
21) Nitrobenzene-d5	8.728	128		61594	34.545	ng/ul	0.00
24) 2-Nitrophenol-d4	9.440	143		75437	35.489	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165		149099	33.494	ng/ul	0.00
31) 4-Chloroaniline-d4	10.481	131		148026	27.873	ng/ul	0.00
46) Dimethylphthalate-d6	13.604	166		554158	36.660	ng/ul	-0.01
49) Acenaphthylene-d8	13.892	160		553118	35.069	ng/ul	0.00
54) 4-Nitrophenol-d4	14.481	143		21910m	9.645	ng/ul	0.03
60) Fluorene-d10	15.210	176		469538	35.429	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.363	200		97157	31.195	ng/ul	-0.01
73) Anthracene-d10	17.116	188		829736	37.692	ng/ul	0.00
81) Pyrene-d10	19.492	212		1066771	37.561	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.451	264		1387505	38.837	ng/ul	-0.01
Target Compounds							
52) Acenaphthene	14.275	153		15098	1.304	ng/ul	Qvalue 96

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

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