

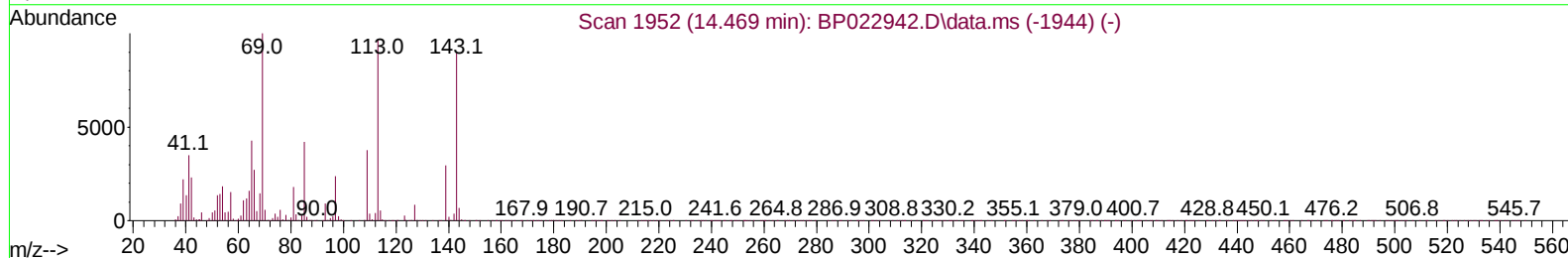
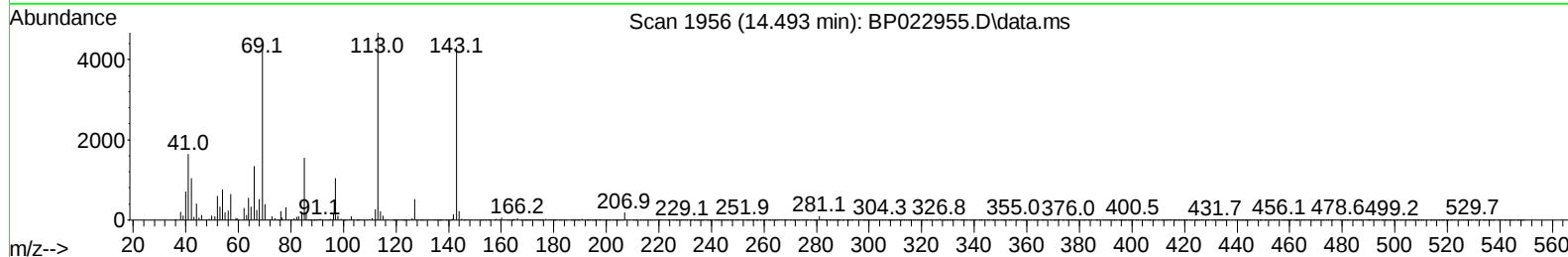
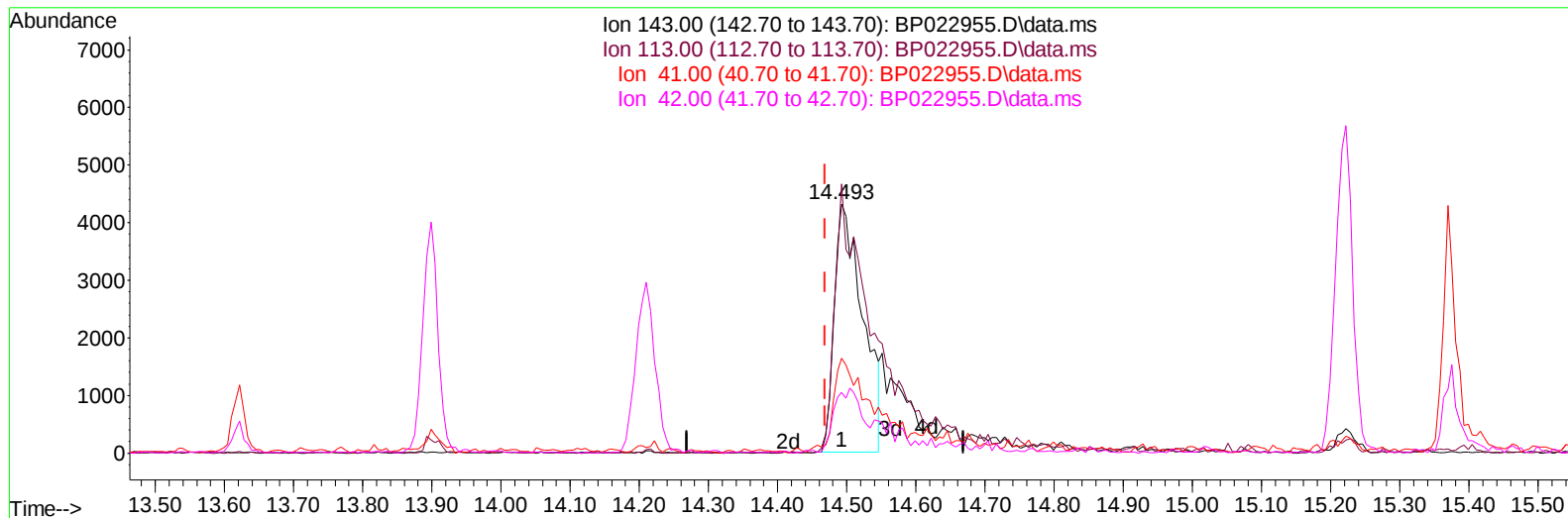
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
Data File : BP022955.D
Acq On : 09 Nov 2024 07:08
Operator : RC/JU
Sample : P4744-13
Misc :
ALS Vial : 40 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP67

Manual IntegrationsAPPROVED

Quant Time: Nov 09 22:14:22 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Nov 08 23:49:48 2024
Response via : Initial Calibration

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



TIC: BP022955.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.493min (+ 0.024) 6.53 ng/u1

response 12273

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	108.05
41.00	39.20	38.00
42.00	25.90	24.23

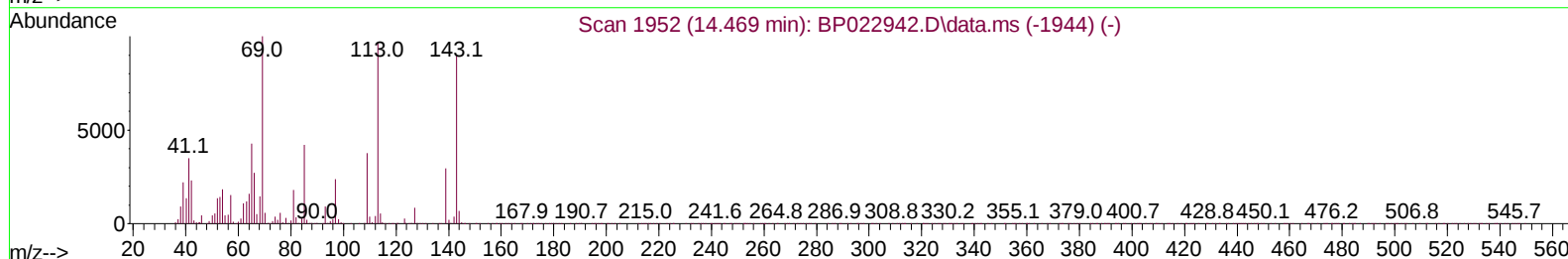
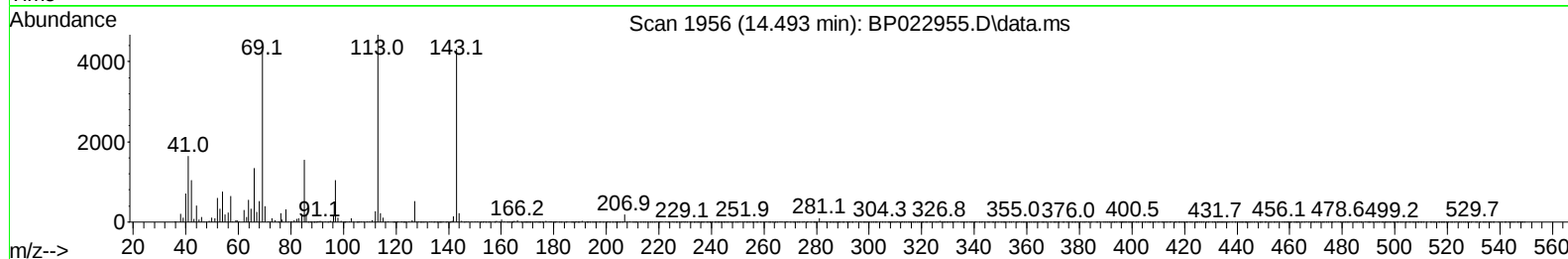
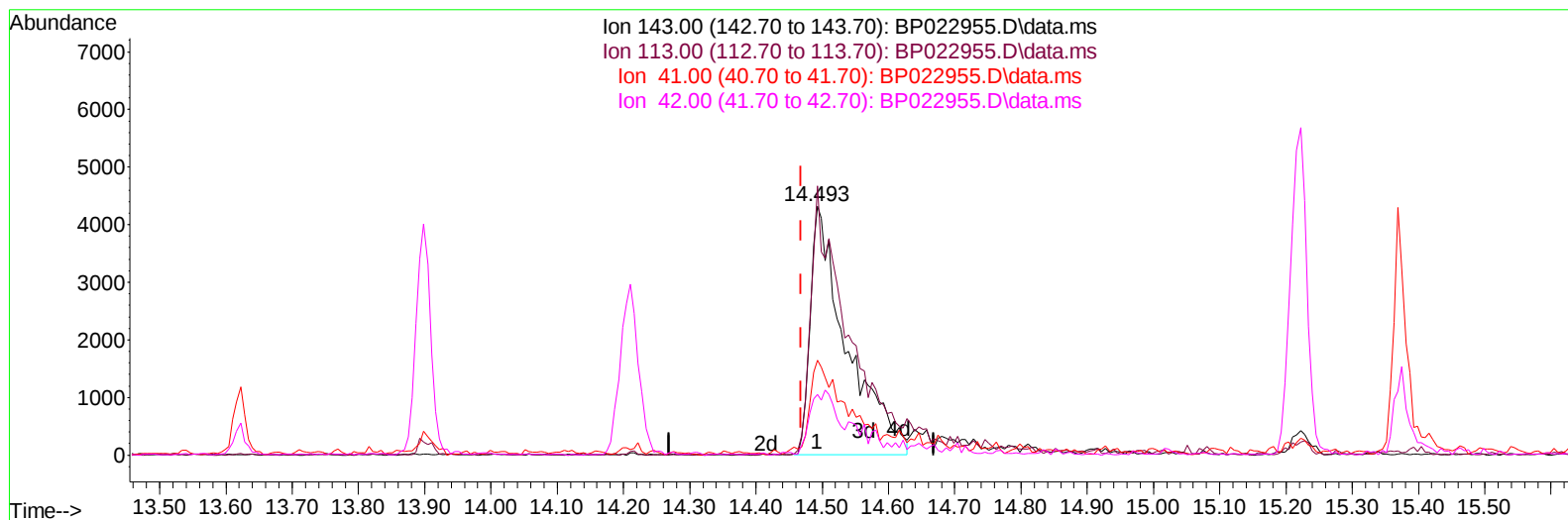
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TIC: BP022955.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.493min (+ 0.024) 8.80 ng/ul m

response 16523

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	108.05
41.00	39.20	38.00
42.00	25.90	24.23

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 Sample : P4744-13
 Misc :
 ALS Vial : 40 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCP67

Manual IntegrationsAPPROVED

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

Quant Time: Nov 09 22:15:02 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP110924.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Nov 08 23:49:48 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	56150	20.000	ng/ul	0.00
20) Naphthalene-d8	10.340	136	213641	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.210	164	173412	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.010	188	440313	20.000	ng/ul	-0.02
79) Chrysene-d12	21.451	240	540731	20.000	ng/ul	0.00
88) Perylene-d12	24.680	264	647335	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	6655	5.697	ng/uL	0.00
4) Pyridine-d5	3.570	84	13276	3.825	ng/ul	0.01
7) Phenol-d5	6.799	99	37735	9.090	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	82258	33.726	ng/ul	0.00
11) 2-Chlorophenol-d4	7.140	132	92532	30.435	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	72687	21.277	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	52395	35.228	ng/ul	0.00
24) 2-Nitrophenol-d4	9.452	143	63151	35.615	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	121201	32.640	ng/ul	0.00
31) 4-Chloroaniline-d4	10.493	131	107829	24.341	ng/ul	0.00
46) Dimethylphthalate-d6	13.622	166	480834	38.472	ng/ul	0.00
49) Acenaphthylene-d8	13.899	160	451891	34.651	ng/ul	0.00
54) 4-Nitrophenol-d4	14.493	143	16523m	8.797	ng/ul	0.02
60) Fluorene-d10	15.222	176	399018	36.414	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.369	200	87397	33.134	ng/ul	0.00
73) Anthracene-d10	17.128	188	723993	38.834	ng/ul	0.00
81) Pyrene-d10	19.498	212	964943	38.743	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.457	264	1220826	39.592	ng/ul	-0.01

Target Compounds Qvalue

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

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