

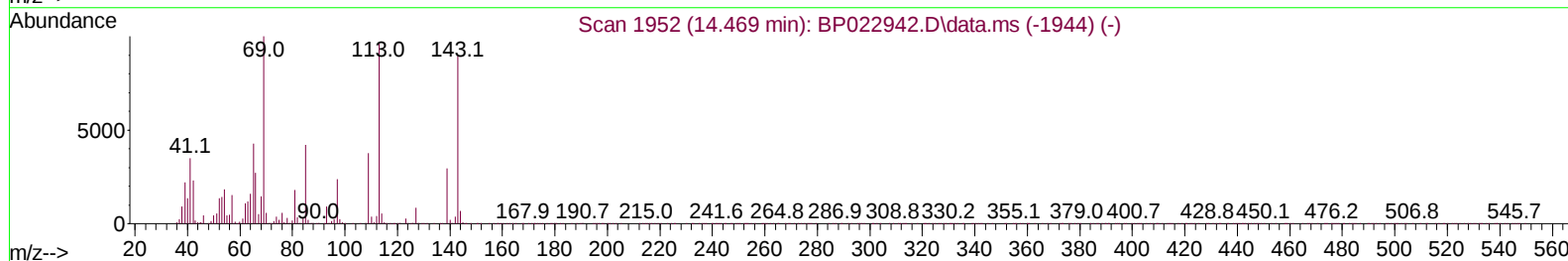
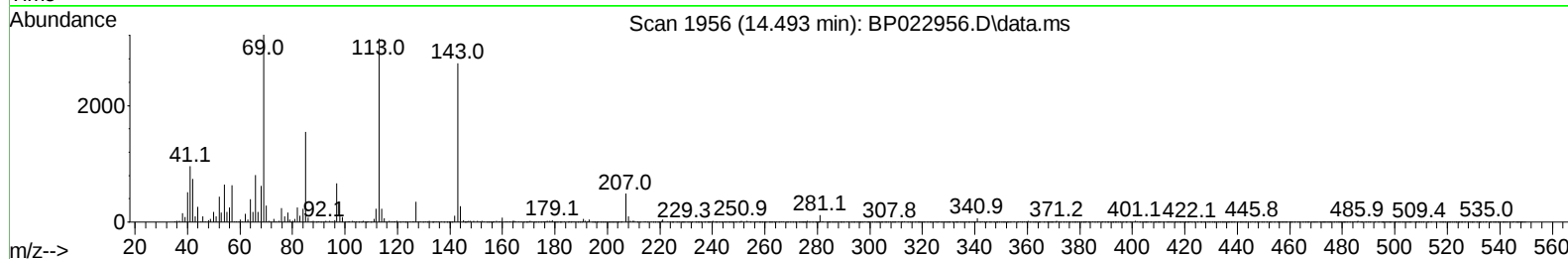
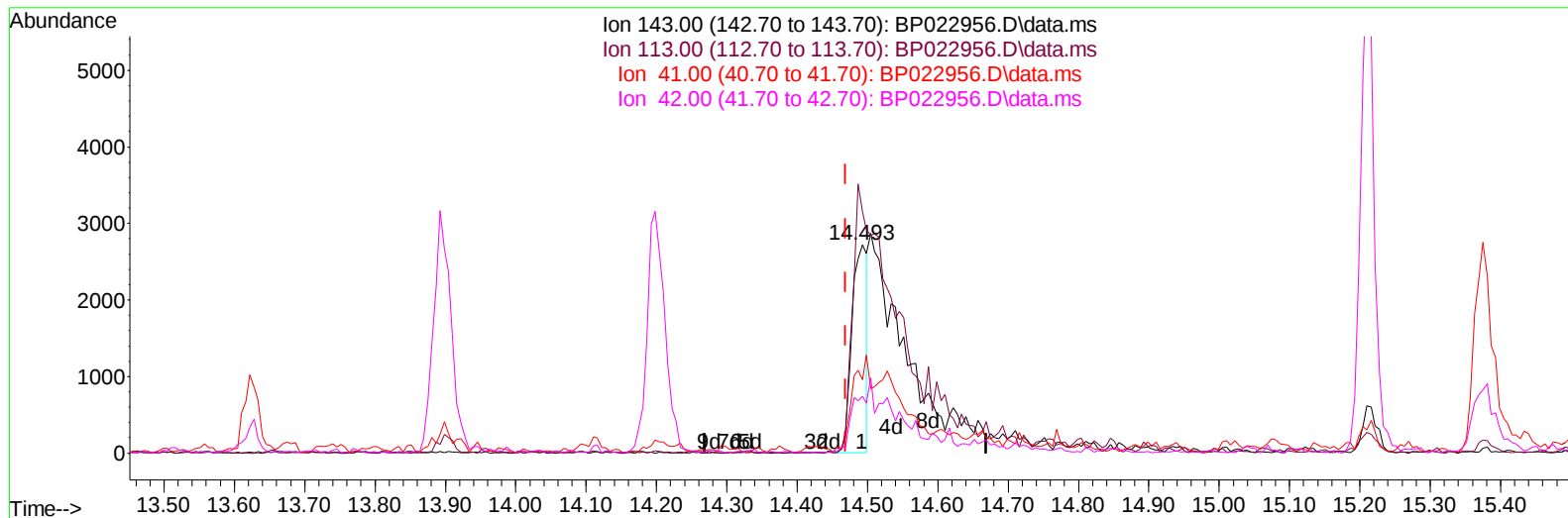
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
Data File : BP022956.D
Acq On : 09 Nov 2024 07:49
Operator : RC/JU
Sample : P4744-14
Misc :
ALS Vial : 41 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP68

Manual IntegrationsAPPROVED

Quant Time: Nov 09 22:16:49 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: BP022956.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.493min (+ 0.024) 2.32 ng/u1

response 4094

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	115.88
41.00	39.20	35.13
42.00	25.90	27.23

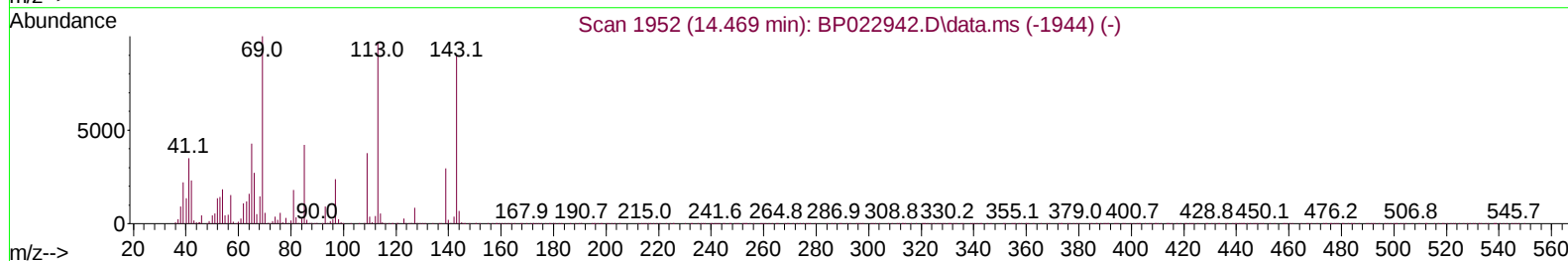
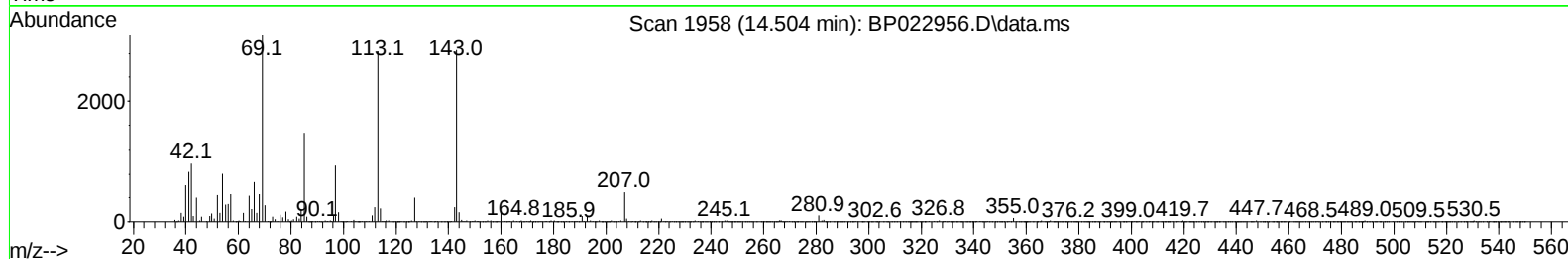
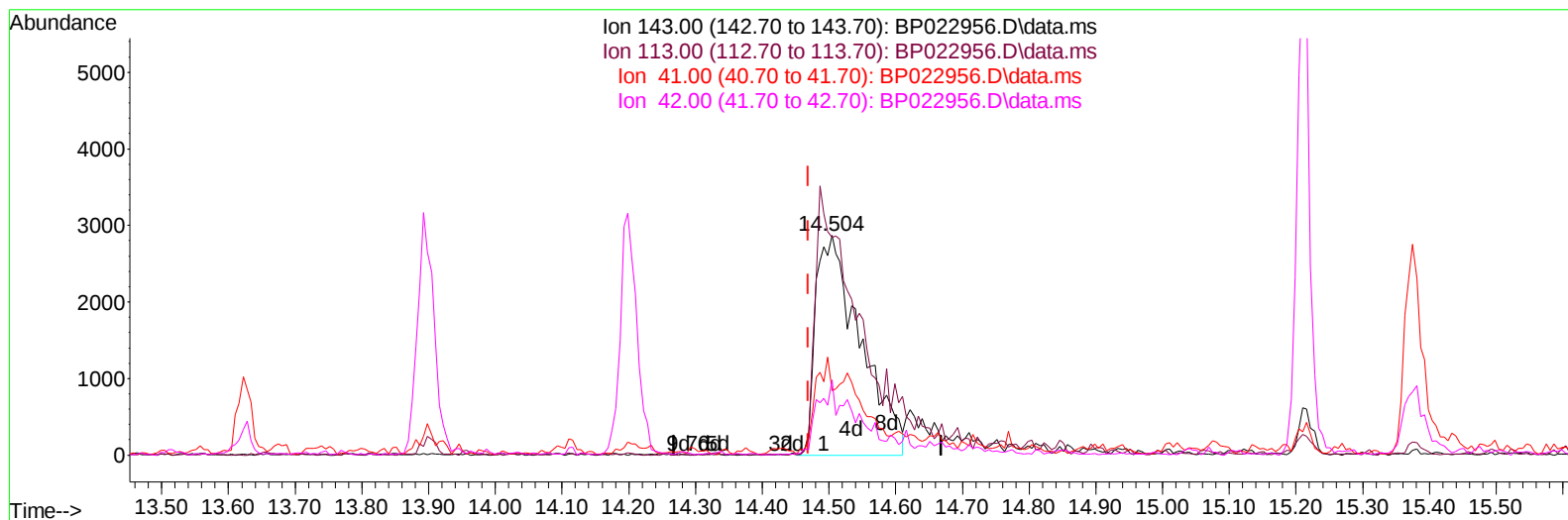
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TIC: BP022956.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.035) 7.55 ng/ul m

response 13351

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	99.23
41.00	39.20	29.39#
42.00	25.90	34.14#

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Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/11/2024
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Quant Time: Nov 09 22:17:36 2024
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 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	55836	20.000	ng/ul	0.00
20) Naphthalene-d8	10.340	136	203509	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.198	164	163205	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.010	188	416609	20.000	ng/ul	-0.02
79) Chrysene-d12	21.451	240	511229	20.000	ng/ul	0.00
88) Perylene-d12	24.668	264	638881	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	5664	4.876	ng/uL	0.00
4) Pyridine-d5	3.564	84	29725	8.613	ng/ul	0.00
7) Phenol-d5	6.799	99	32974	7.988	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	71429	29.451	ng/ul	0.00
11) 2-Chlorophenol-d4	7.140	132	82622	27.328	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	62647	18.441	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	44231	31.219	ng/ul	0.00
24) 2-Nitrophenol-d4	9.452	143	54071	32.013	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	104120	29.436	ng/ul	0.00
31) 4-Chloroaniline-d4	10.493	131	114367	27.102	ng/ul	0.00
46) Dimethylphthalate-d6	13.622	166	413763	35.176	ng/ul	0.00
49) Acenaphthylene-d8	13.898	160	390905	31.849	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	13351m	7.553	ng/ul	0.04
60) Fluorene-d10	15.210	176	356098	34.529	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.375	200	74213	29.736	ng/ul	0.00
73) Anthracene-d10	17.122	188	655954	37.186	ng/ul	0.00
81) Pyrene-d10	19.504	212	867759	36.851	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.457	264	1130009	37.132	ng/ul	-0.01

Target Compounds Qvalue

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

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