

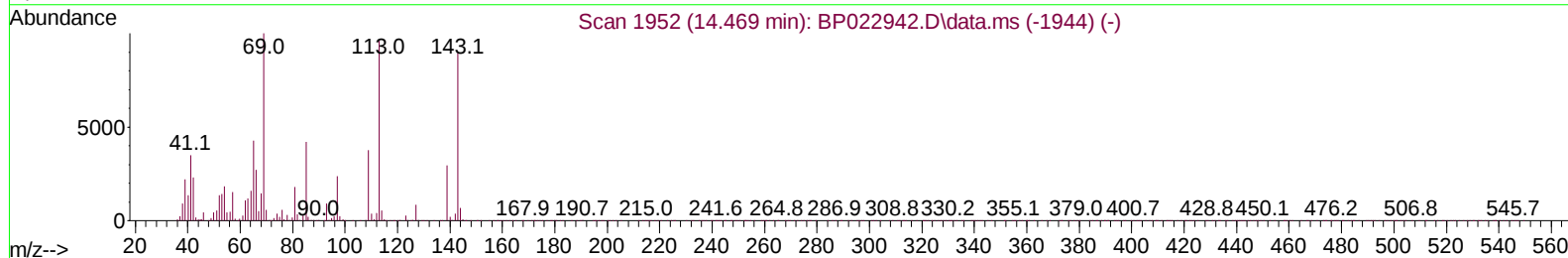
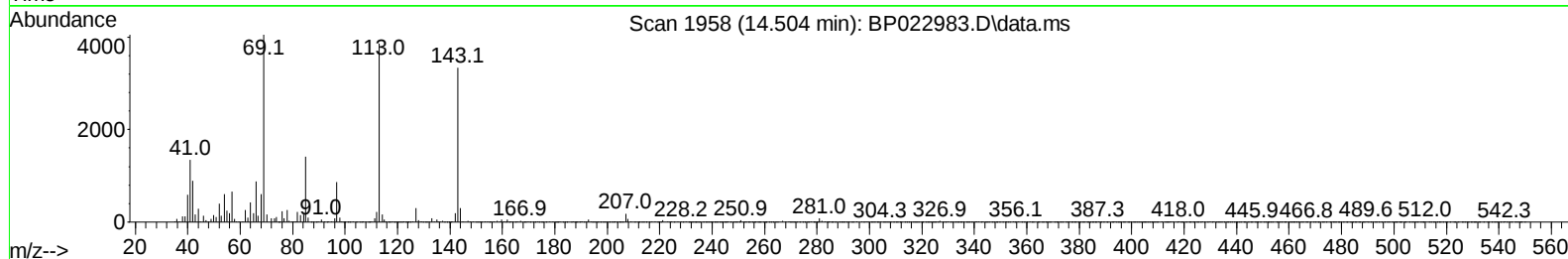
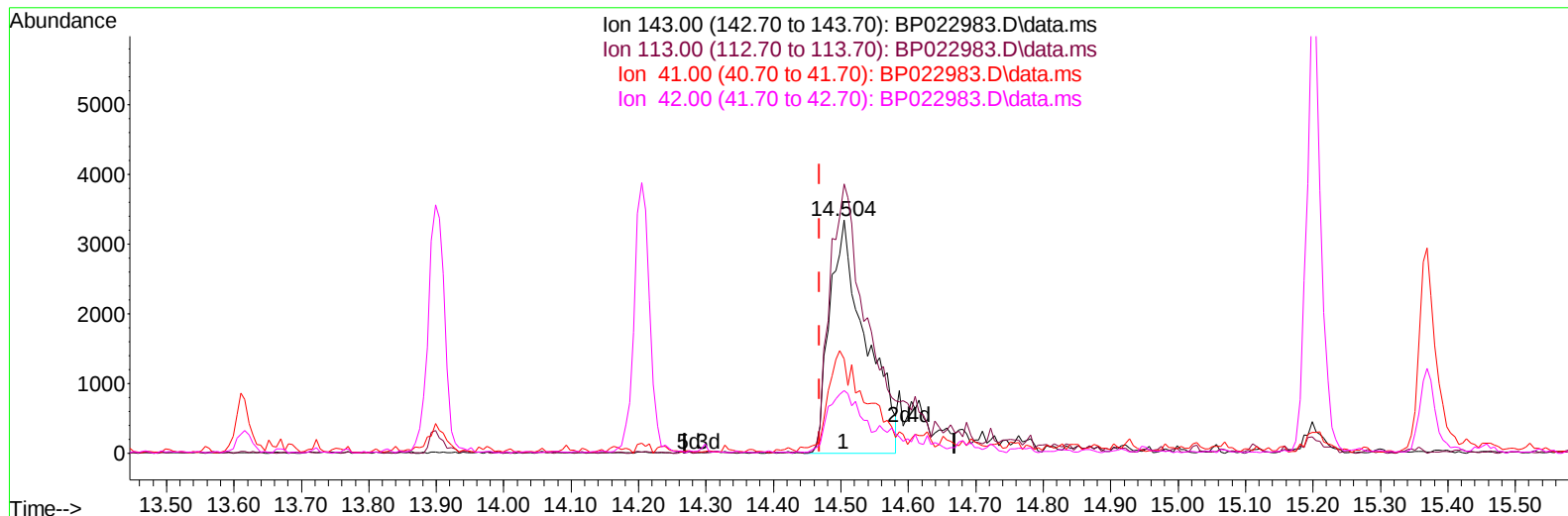
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
Data File : BP022983.D
Acq On : 10 Nov 2024 02:45
Operator : RC/JU
Sample : P4746-13
Misc :
ALS Vial : 70 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP86

Manual IntegrationsAPPROVED

Quant Time: Nov 10 03:17:53 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: BP022983.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.036) 6.33 ng/u1

response 12297

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	115.52
41.00	39.20	40.35
42.00	25.90	26.74

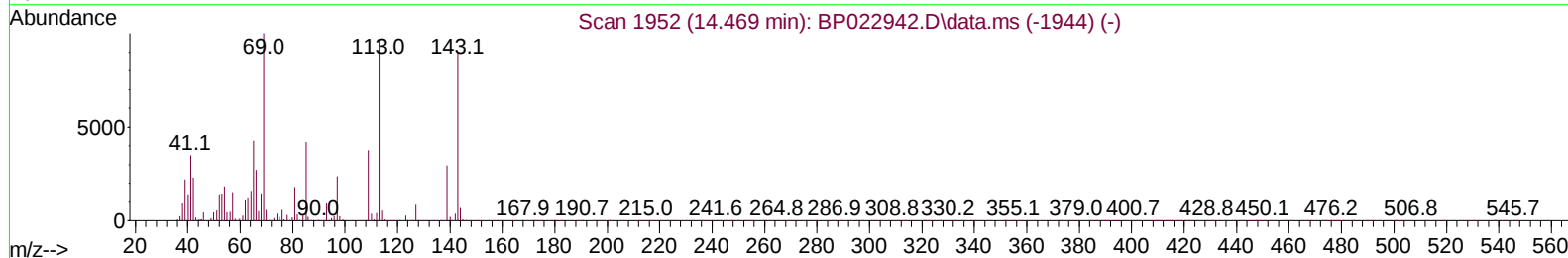
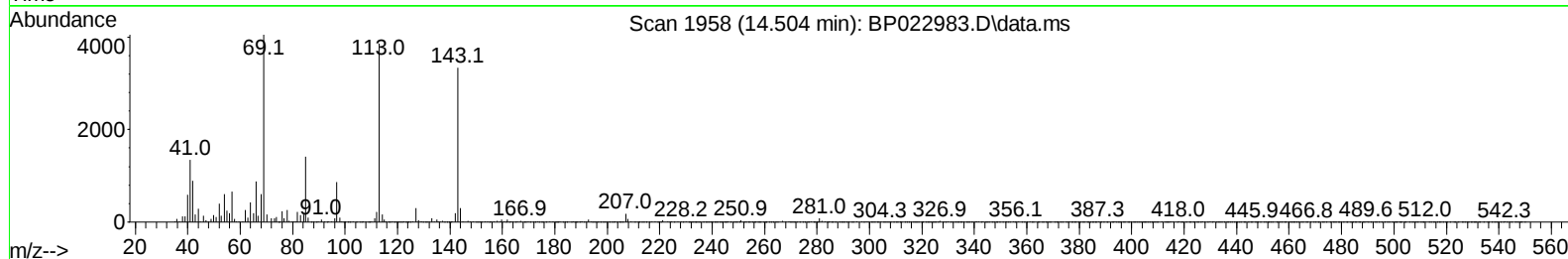
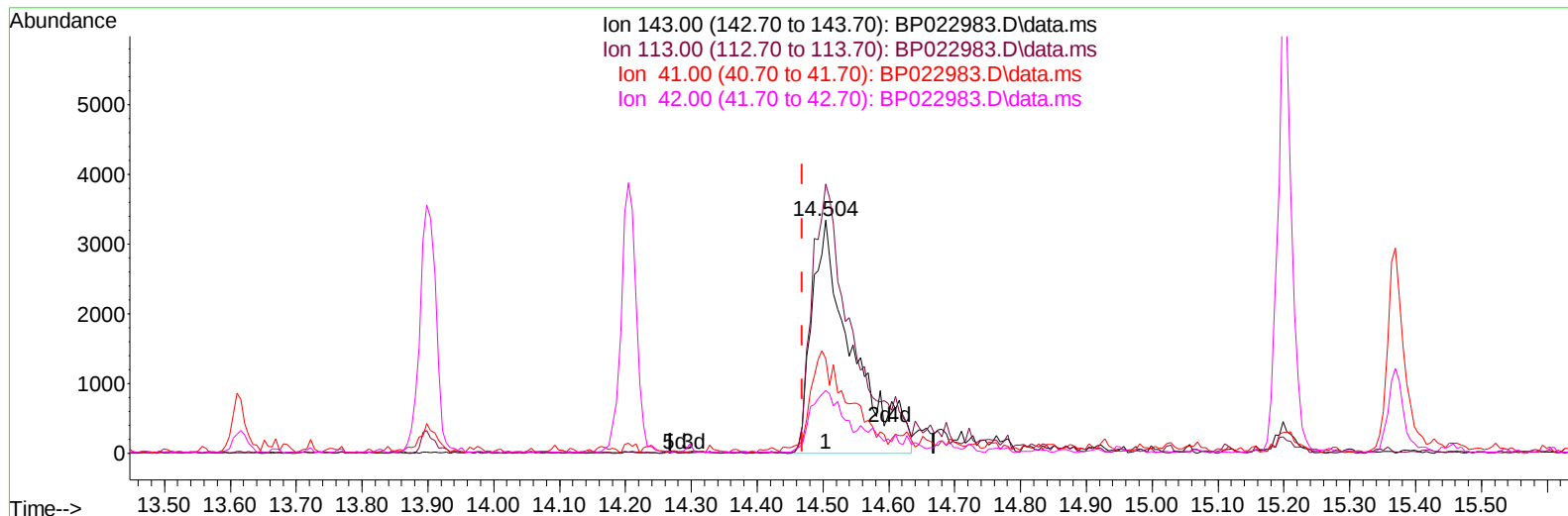
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TIC: BP022983.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.036) 7.27 ng/ul m

response 14110

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	115.52
41.00	39.20	40.35
42.00	25.90	26.74

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Reviewed By :Yogesh Patel 11/11/2024
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Quant Time: Nov 10 03:18:51 2024
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 QLast Update : Fri Nov 08 23:49:48 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	56486	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.340	136	217784	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.204	164	179271	20.000	ng/ul	-0.02
64) Phenanthrene-d10	16.998	188	470040	20.000	ng/ul	-0.04
79) Chrysene-d12	21.451	240	567008	20.000	ng/ul	0.00
88) Perylene-d12	24.669	264	693759	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	5113	4.351	ng/uL	0.00
4) Pyridine-d5	3.558	84	31091	8.905	ng/ul	0.00
7) Phenol-d5	6.793	99	34225	8.195	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.934	67	72388	29.503	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.140	132	80832	26.428	ng/ul	0.00
15) 4-Methylphenol-d8	8.305	113	65374	19.023	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	46050	30.373	ng/ul	0.00
24) 2-Nitrophenol-d4	9.446	143	55861	30.905	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	104947	27.725	ng/ul	0.00
31) 4-Chloroaniline-d4	10.487	131	121884	26.990	ng/ul	-0.01
46) Dimethylphthalate-d6	13.616	166	427524	33.089	ng/ul	0.00
49) Acenaphthylene-d8	13.898	160	415882	30.848	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	14110m	7.267	ng/ul	0.04
60) Fluorene-d10	15.198	176	369711	32.637	ng/ul	-0.04
65) 4,6-Dinitro-2-methylph...	15.369	200	78156	27.757	ng/ul	0.00
73) Anthracene-d10	17.116	188	665630	33.445	ng/ul	0.00
81) Pyrene-d10	19.498	212	889884	34.073	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.439	264	1110513	33.605	ng/ul	-0.03

Target Compounds Qvalue

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

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