

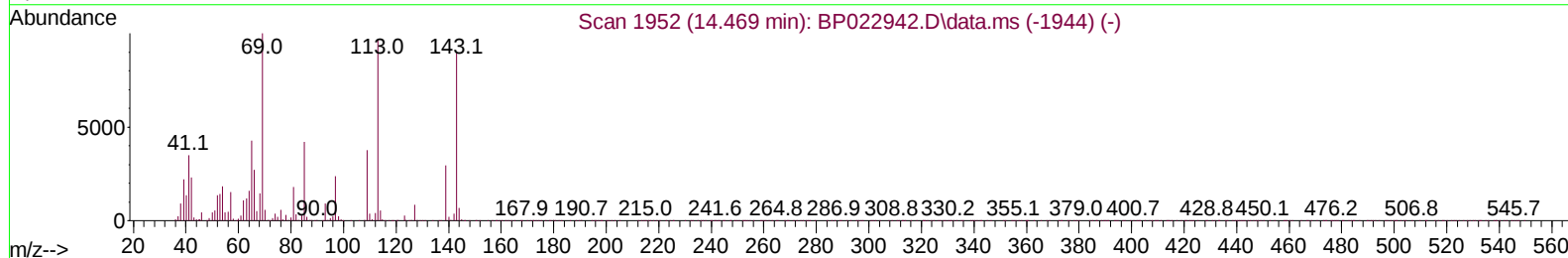
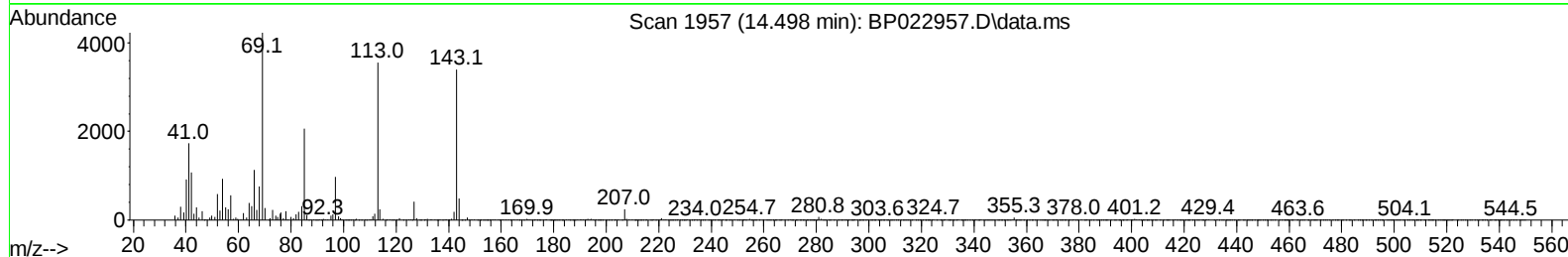
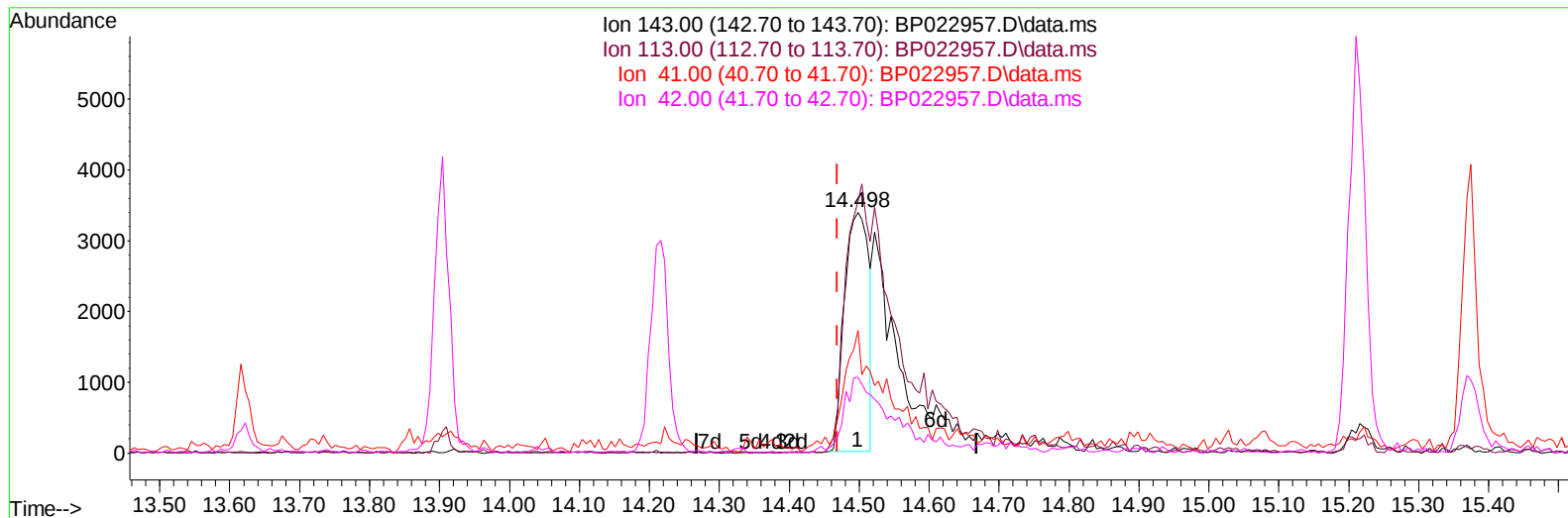
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
 Data File : BP022957.D
 Acq On : 09 Nov 2024 08:29
 Operator : RC/JU
 Sample : P4744-15
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCP70

Manual IntegrationsAPPROVED

Quant Time: Nov 09 22:18:58 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: BP022957.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.030) 4.59 ng/u1

response 8218

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	104.58
41.00	39.20	50.90#
42.00	25.90	31.56#

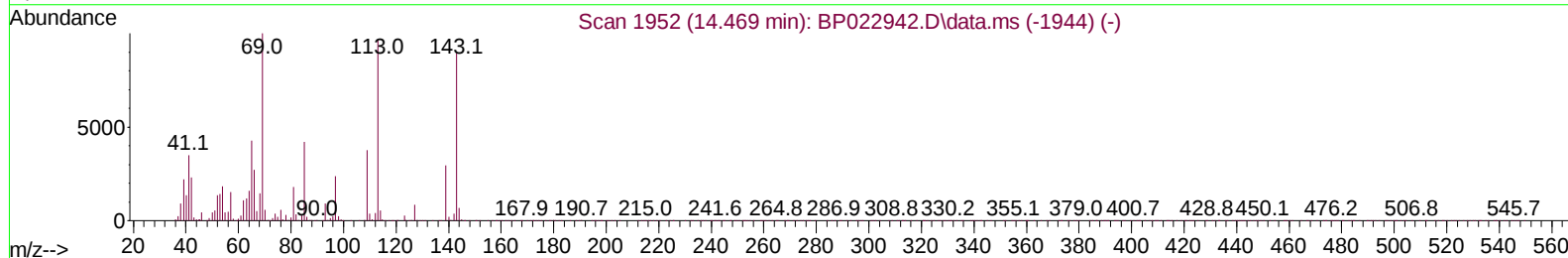
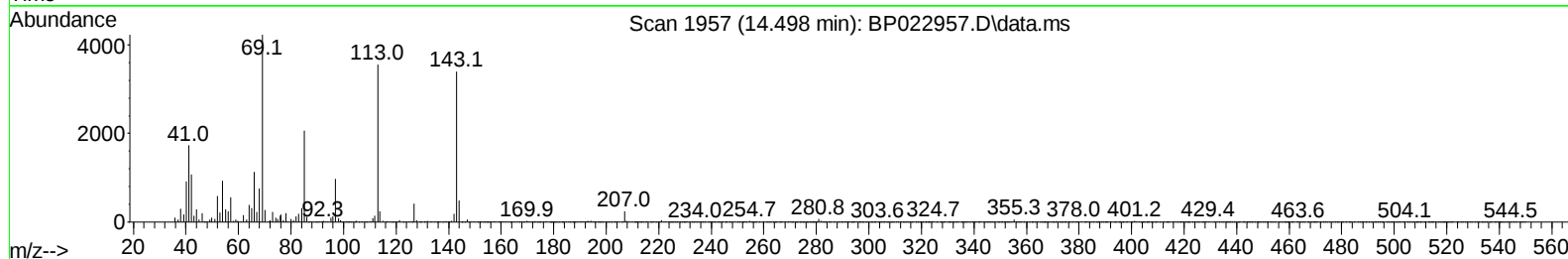
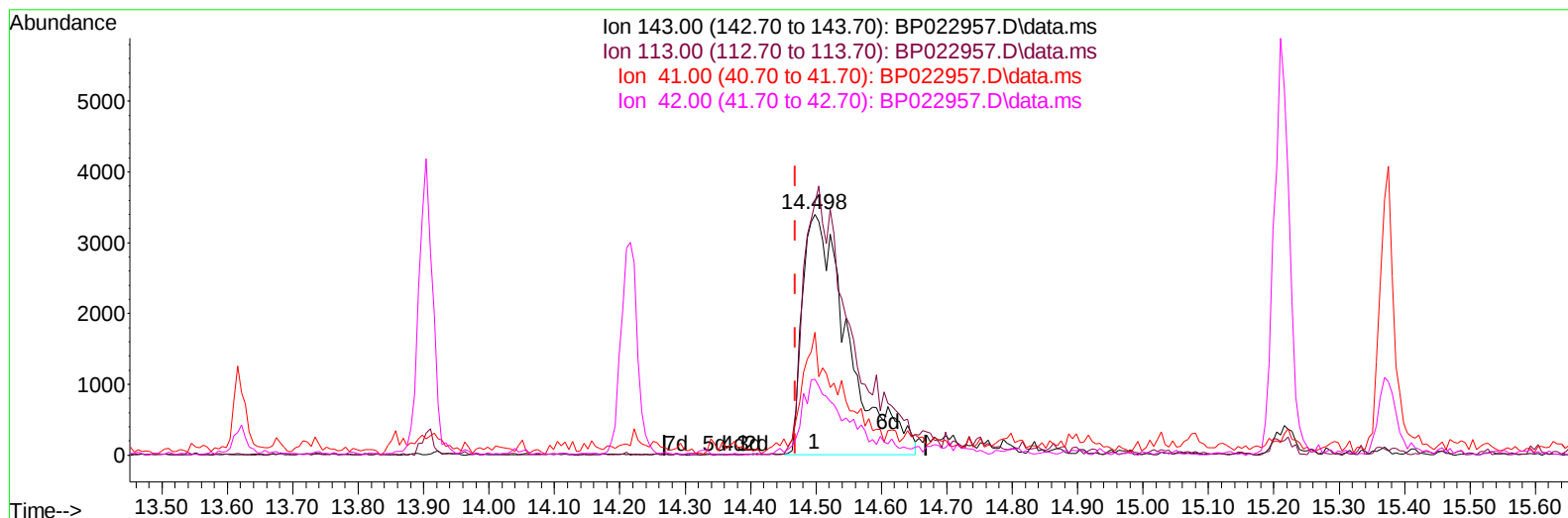
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TIC: BP022957.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (+ 0.030) 9.27 ng/ul m

response 16596

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	104.58
41.00	39.20	50.90#
42.00	25.90	31.56#

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Quant Time: Nov 09 22:19:43 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	55105	20.000	ng/ul	0.00
20) Naphthalene-d8	10.346	136	204750	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.216	164	165270	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.016	188	415008	20.000	ng/ul	-0.02
79) Chrysene-d12	21.445	240	544772	20.000	ng/ul	-0.01
88) Perylene-d12	24.674	264	678511	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	6400	5.583	ng/uL	0.00
4) Pyridine-d5	3.569	84	18894	5.547	ng/ul	0.01
7) Phenol-d5	6.805	99	38392	9.424	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	6.946	67	80518	33.639	ng/ul	0.00
11) 2-Chlorophenol-d4	7.140	132	91608	30.702	ng/ul	0.00
15) 4-Methylphenol-d8	8.310	113	75128	22.409	ng/ul	0.00
21) Nitrobenzene-d5	8.740	128	51024	35.796	ng/ul	0.00
24) 2-Nitrophenol-d4	9.452	143	61002	35.897	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.993	165	119145	33.479	ng/ul	0.00
31) 4-Chloroaniline-d4	10.499	131	81163	19.117	ng/ul	0.00
46) Dimethylphthalate-d6	13.622	166	452134	37.958	ng/ul	0.00
49) Acenaphthylene-d8	13.904	160	443054	35.647	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	16596m	9.271	ng/ul	0.03
60) Fluorene-d10	15.210	176	386256	36.986	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.375	200	82356	33.127	ng/ul	0.00
73) Anthracene-d10	17.128	188	671296	38.203	ng/ul	0.00
81) Pyrene-d10	19.492	212	962085	38.341	ng/ul	-0.02
92) Benzo(a)pyrene-d12	24.439	264	1261514	39.032	ng/ul	-0.03

Target Compounds Qvalue

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

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