

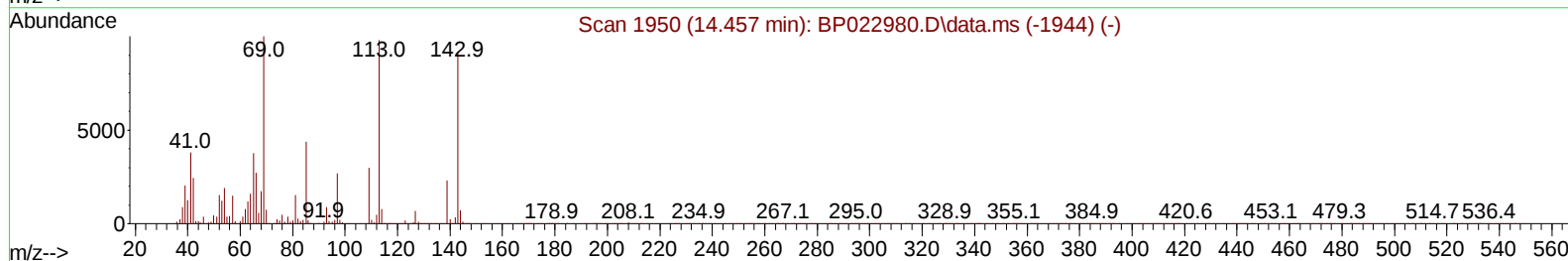
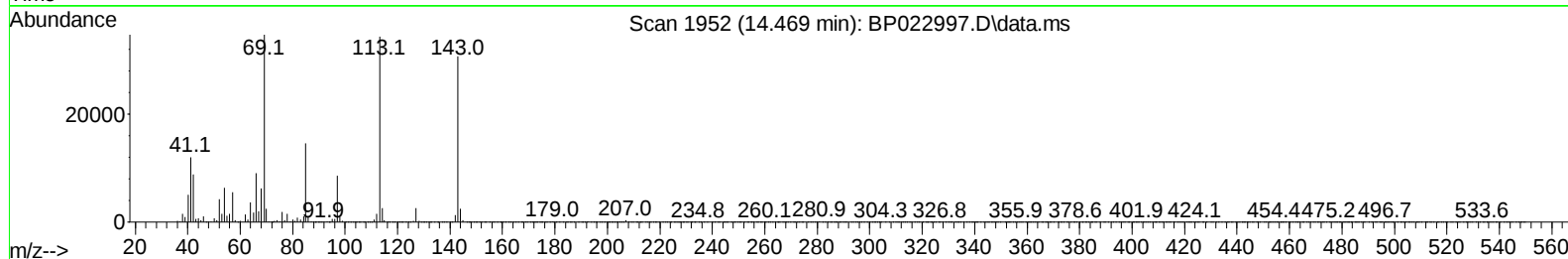
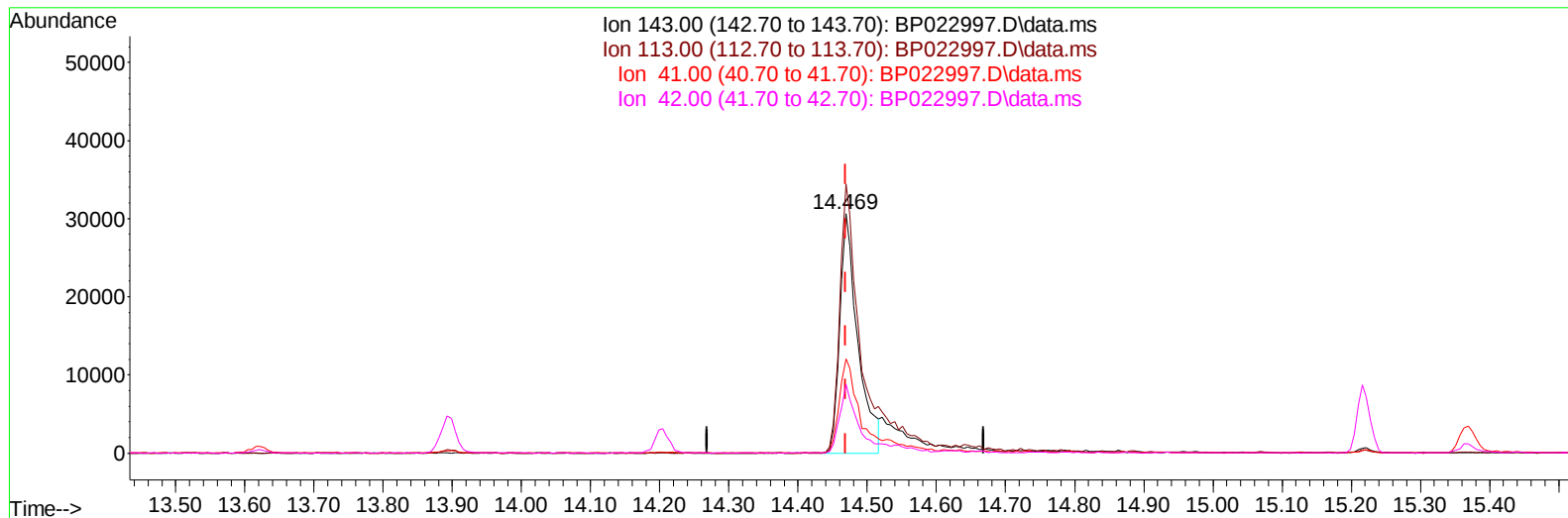
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
Data File : BP022997.D
Acq On : 10 Nov 2024 12:53
Operator : RC/JU
Sample : P4746-22
Misc :
ALS Vial : 85 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP95

Manual IntegrationsAPPROVED

Quant Time: Nov 10 22:39:40 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: BP022997.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.469min (+ 0.000) 36.73 ng/u1

response 55684

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	112.21
41.00	39.20	39.29
42.00	25.90	28.75

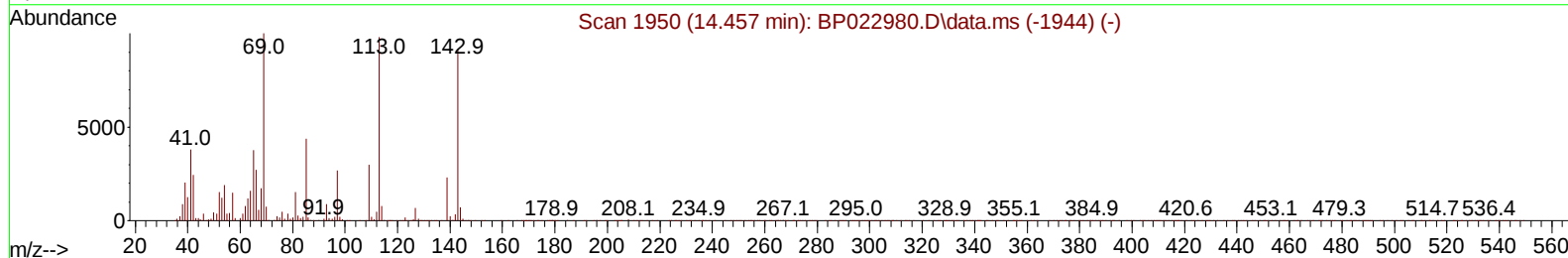
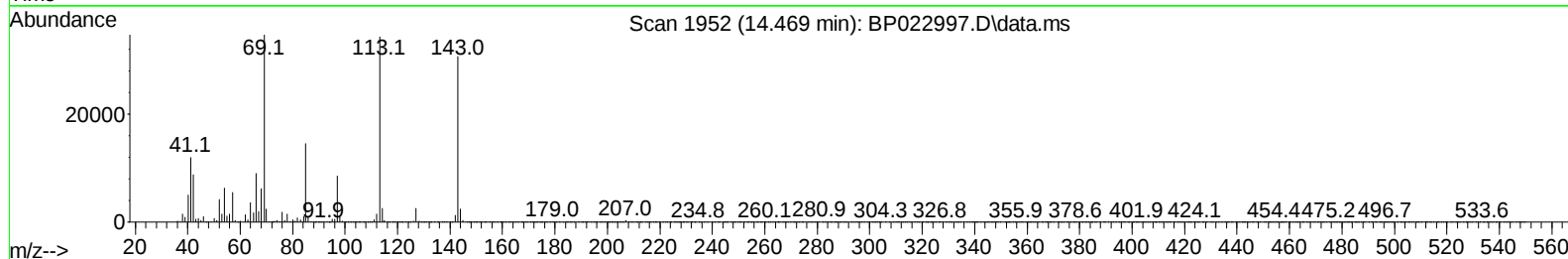
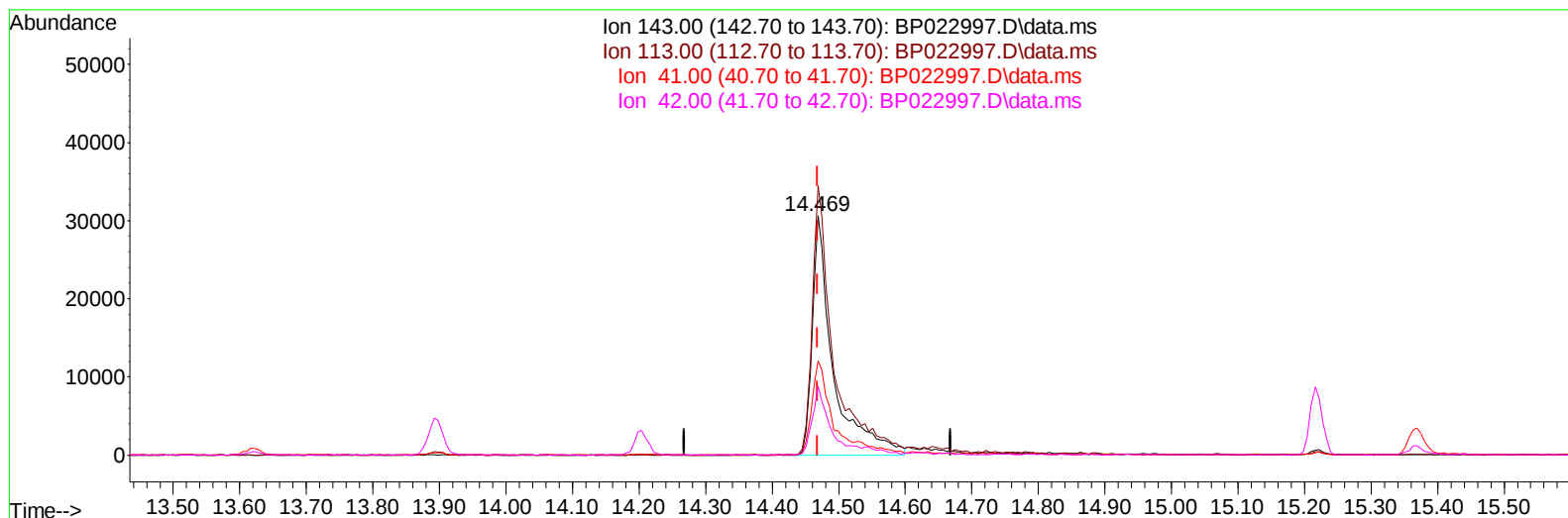
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(54) 4-Nitrophenol-d4 (S)

14.469min (+ 0.000) 44.36 ng/ul m

response 67257

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	112.21
41.00	39.20	39.29
42.00	25.90	28.75

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Reviewed By :Yogesh Patel 11/11/2024
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Quant Time: Nov 10 23:05:55 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	7.587	152	42970	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.334	136	166126	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.204	164	139970	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.010	188	351499	20.000	ng/ul	-0.02
79) Chrysene-d12	21.445	240	413160	20.000	ng/ul	#-0.01
88) Perylene-d12	24.674	264	476586	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	7696	8.609	ng/uL	0.00
4) Pyridine-d5	3.552	84	106593	40.133	ng/ul	0.00
7) Phenol-d5	6.793	99	140142	44.113	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	84269	45.148	ng/ul	0.00
11) 2-Chlorophenol-d4	7.140	132	107843	46.351	ng/ul	0.00
15) 4-Methylphenol-d8	8.299	113	119394	45.670	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	53745	46.471	ng/ul	0.00
24) 2-Nitrophenol-d4	9.446	143	67768	49.151	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.981	165	132931	46.038	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.493	131	156442	45.415	ng/ul	0.00
46) Dimethylphthalate-d6	13.622	166	521966	51.741	ng/ul	0.00
49) Acenaphthylene-d8	13.892	160	506689	48.136	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.469	143	67257m	44.363	ng/ul	0.00
60) Fluorene-d10	15.216	176	441963	49.969	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.369	200	89024	42.279	ng/ul	0.00
73) Anthracene-d10	17.104	188	793435	53.312	ng/ul	-0.02
81) Pyrene-d10	19.498	212	1026101	53.919	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.445	264	1238299	54.547	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.175	88	4695	4.545	ng/uL#	93

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

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