

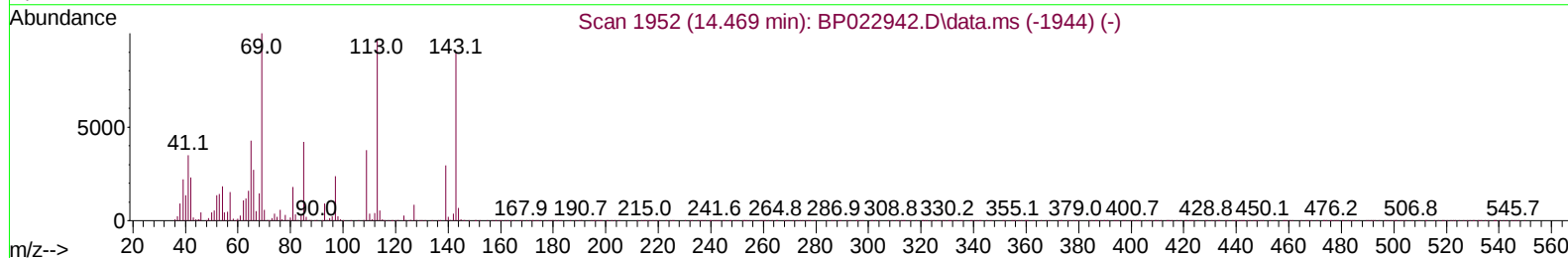
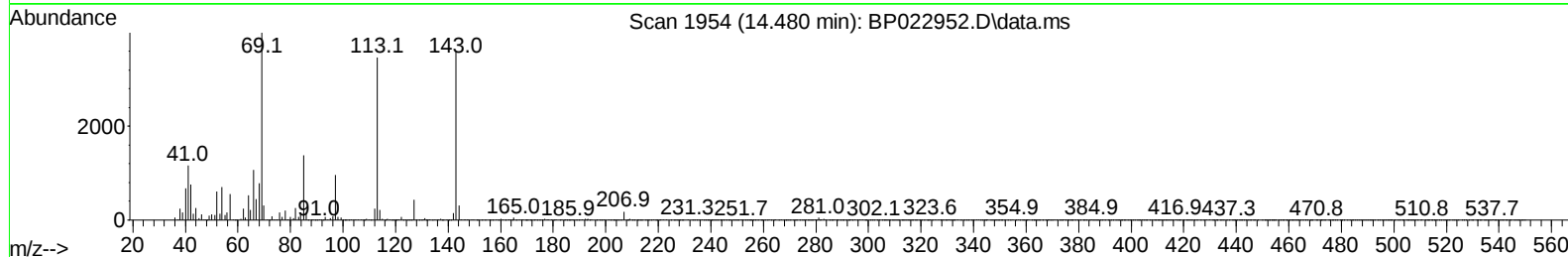
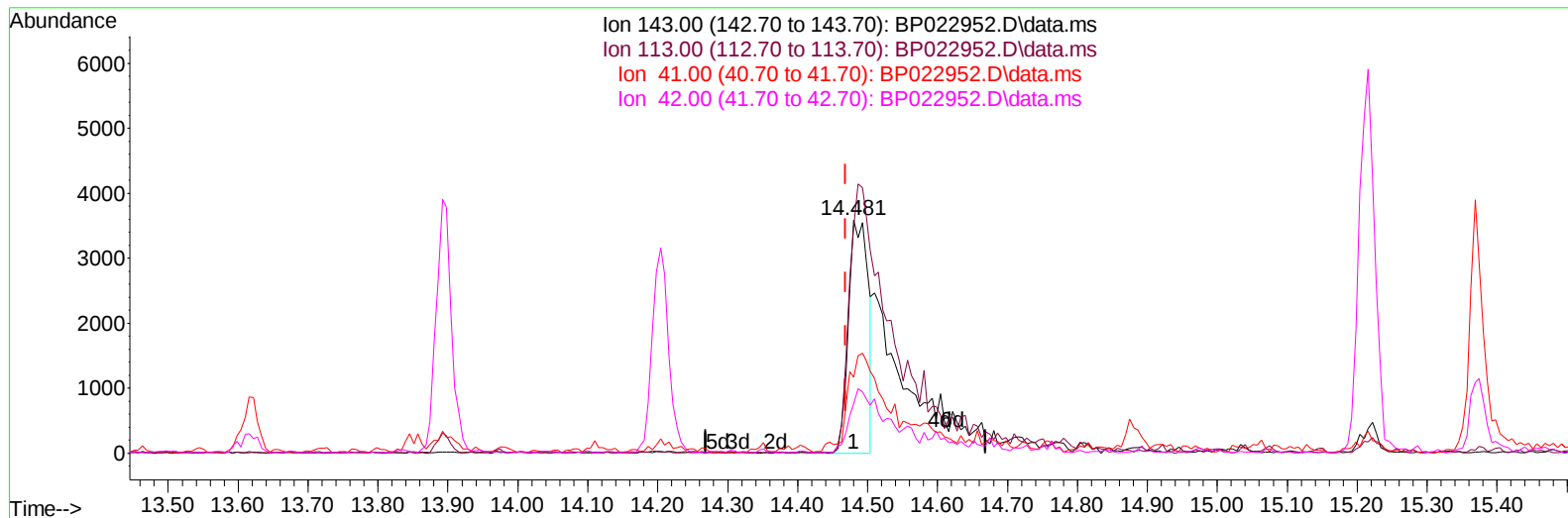
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110924\
Data File : BP022952.D
Acq On : 09 Nov 2024 05:07
Operator : RC/JU
Sample : P4744-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP64

Manual IntegrationsAPPROVED

Quant Time: Nov 09 05:34:05 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: BP022952.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.480min (+ 0.012) 4.50 ng/u1

response 7088

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	96.60
41.00	39.20	32.43
42.00	25.90	21.04

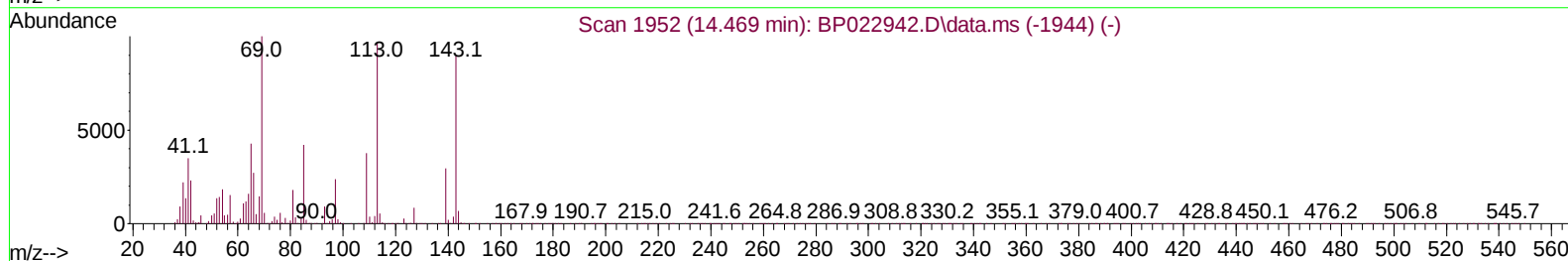
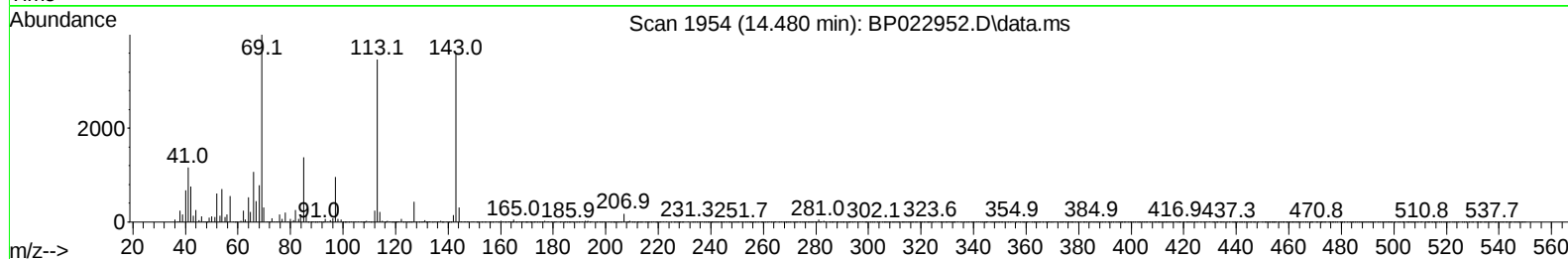
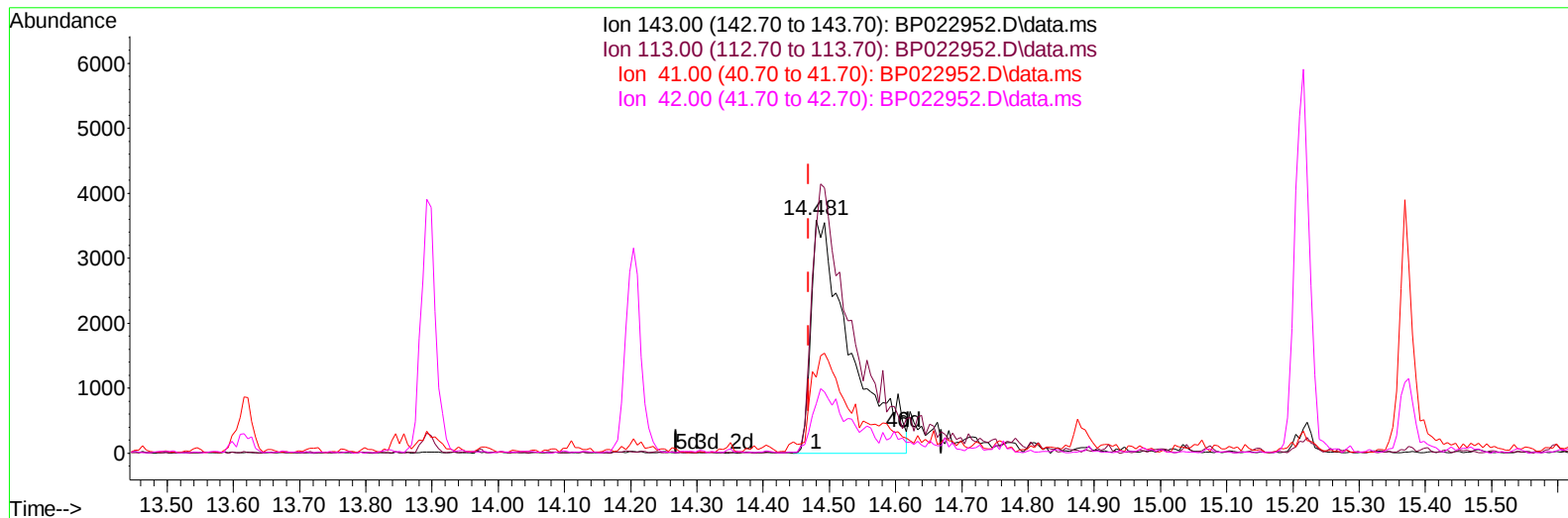
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(54) 4-Nitrophenol-d4 (S)

14.480min (+ 0.012) 9.36 ng/ul m

response 14750

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	107.90	96.60
41.00	39.20	32.43
42.00	25.90	21.04

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.587	152	47277	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.340	136	180149	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.204	164	145437	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.010	188	359831	20.000	ng/ul	-0.02
79) Chrysene-d12	21.457	240	484650	20.000	ng/ul	0.00
88) Perylene-d12	24.680	264	615688	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	5946	6.045	ng/uL	0.00
4) Pyridine-d5	3.563	84	20356	6.966	ng/ul	0.00
7) Phenol-d5	6.799	99	35219	10.076	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	71663	34.897	ng/ul	0.00
11) 2-Chlorophenol-d4	7.140	132	81471	31.826	ng/ul	0.00
15) 4-Methylphenol-d8	8.298	113	67826	23.581	ng/ul	0.00
21) Nitrobenzene-d5	8.734	128	46041	36.711	ng/ul	0.00
24) 2-Nitrophenol-d4	9.445	143	55933	37.409	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.987	165	101260	32.339	ng/ul	0.00
31) 4-Chloroaniline-d4	10.498	131	77580	20.768	ng/ul	0.00
46) Dimethylphthalate-d6	13.616	166	420028	40.071	ng/ul	0.00
49) Acenaphthylene-d8	13.898	160	397705	36.362	ng/ul	0.00
54) 4-Nitrophenol-d4	14.480	143	14750m	9.364	ng/ul	0.01
60) Fluorene-d10	15.216	176	352858	38.395	ng/ul	-0.02
65) 4,6-Dinitro-2-methylph...	15.375	200	78095	36.229	ng/ul	0.00
73) Anthracene-d10	17.122	188	636423	41.772	ng/ul	0.00
81) Pyrene-d10	19.498	212	880543	39.445	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.451	264	1192903	40.675	ng/ul	-0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.175	88	1420	1.249	ng/uL	91

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

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