

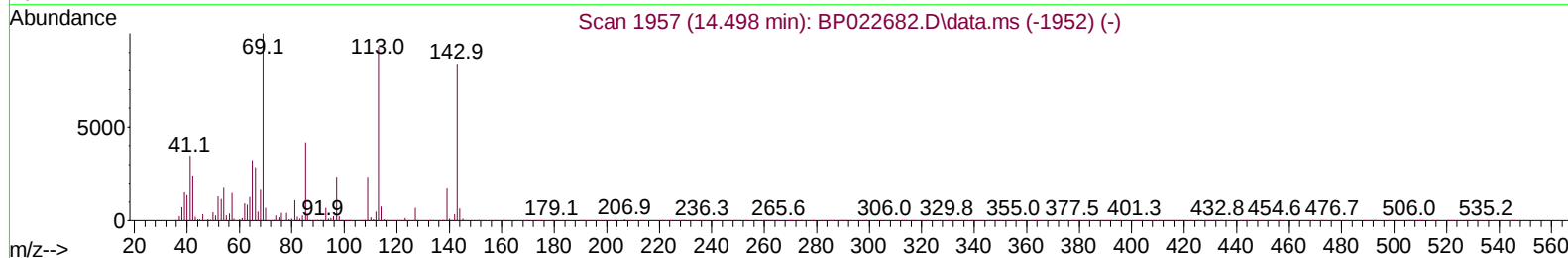
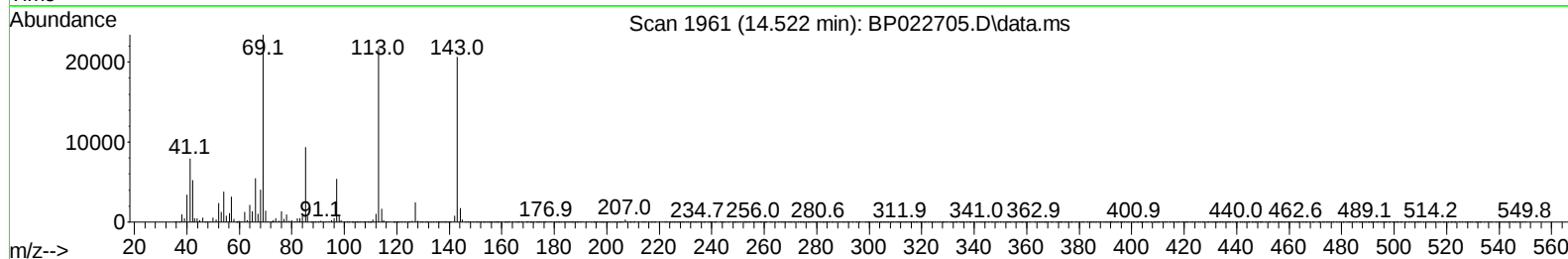
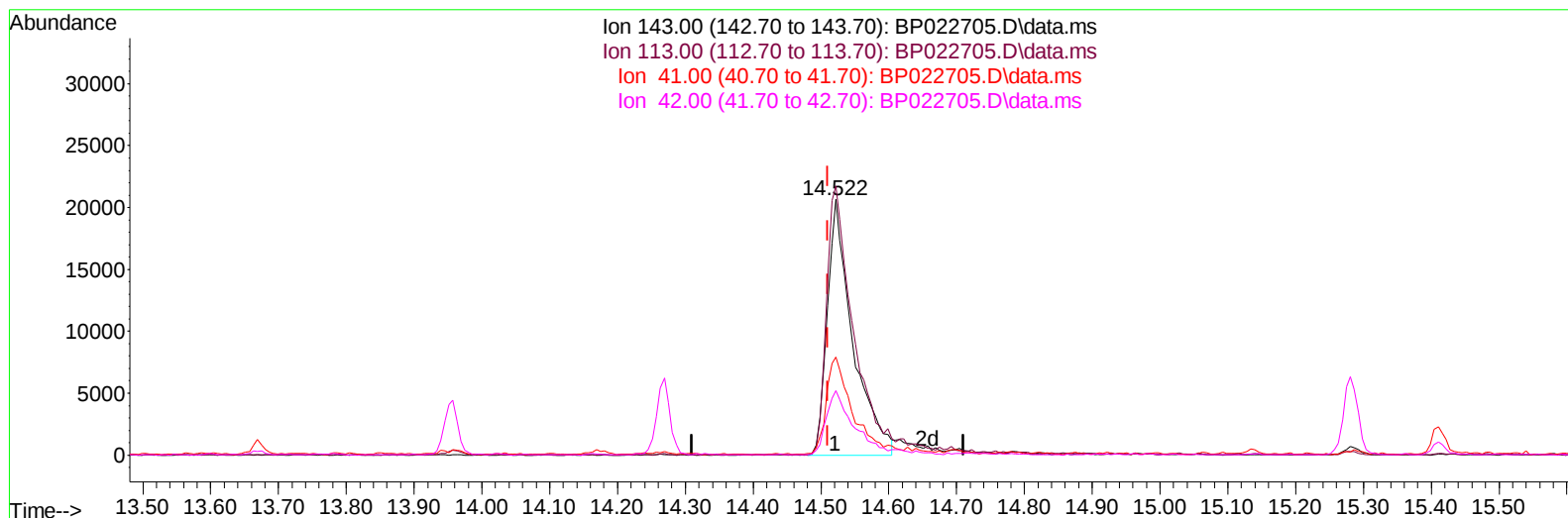
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP102824\
Data File : BP022705.D
Acq On : 29 Oct 2024 08:25
Operator : RC/JU
Sample : P4530-16
Misc :
ALS Vial : 33 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
F7H43

Manual IntegrationsAPPROVED

Quant Time: Oct 29 08:53:59 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Oct 29 04:53:07 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 10/30/2024
Supervised By :mohammad ahmed 10/30/2024



TIC: BP022705.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.012) 14.35 ng/u1

response 53682

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	105.23
41.00	35.90	38.46
42.00	23.70	25.32

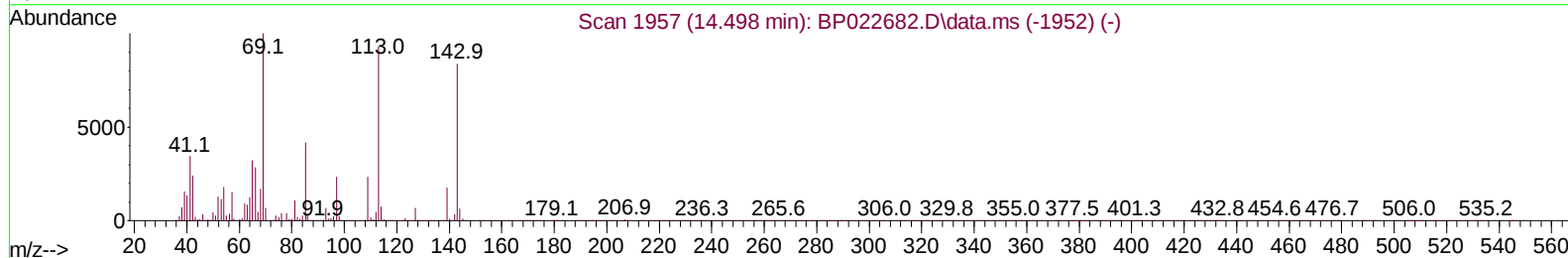
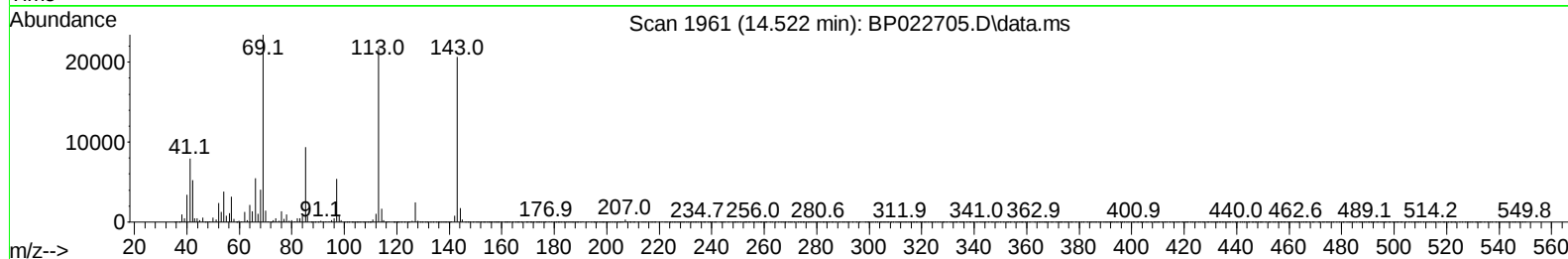
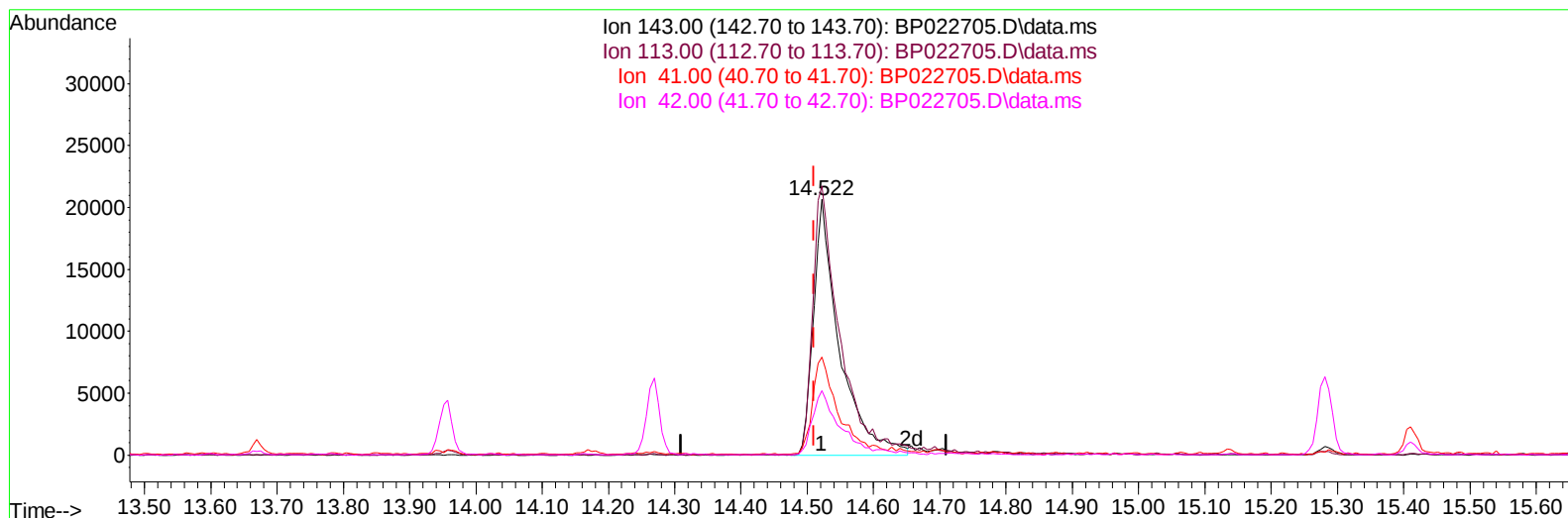
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TIC: BP022705.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.012) 15.03 ng/ul m

response 56222

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	105.23
41.00	35.90	38.46
42.00	23.70	25.32

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.646	152	89531	20.000	ng/ul	0.00
20) Naphthalene-d8	10.399	136	351050	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.269	164	280644	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.069	188	685562	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	703728	20.000	ng/ul	# 0.00
88) Perylene-d12	24.721	264	843275	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.187	96	7577	3.289	ng/uL	0.00
4) Pyridine-d5	3.599	84	78569	12.422	ng/ul	0.00
7) Phenol-d5	6.840	99	147832	19.616	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.987	67	86519	19.532	ng/ul	0.00
11) 2-Chlorophenol-d4	7.187	132	110396	20.644	ng/ul	0.00
15) 4-Methylphenol-d8	8.352	113	124266	20.661	ng/ul	0.00
21) Nitrobenzene-d5	8.787	128	55844	20.689	ng/ul	0.00
24) 2-Nitrophenol-d4	9.499	143	65800	21.056	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.040	165	137056	20.635	ng/ul	0.00
31) 4-Chloroaniline-d4	10.546	131	139603	17.788	ng/ul	0.00
46) Dimethylphthalate-d6	13.669	166	479065	21.485	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	515668	21.774	ng/ul	0.01
54) 4-Nitrophenol-d4	14.522	143	56222m	15.030	ng/ul	0.01
60) Fluorene-d10	15.281	176	423747	21.910	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.410	200	52717	11.692	ng/ul	0.00
73) Anthracene-d10	17.157	188	703143	21.803	ng/ul	0.00
81) Pyrene-d10	19.545	212	863533	23.204	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.516	264	1028451	22.837	ng/ul	0.00

Target Compounds Qvalue

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

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