

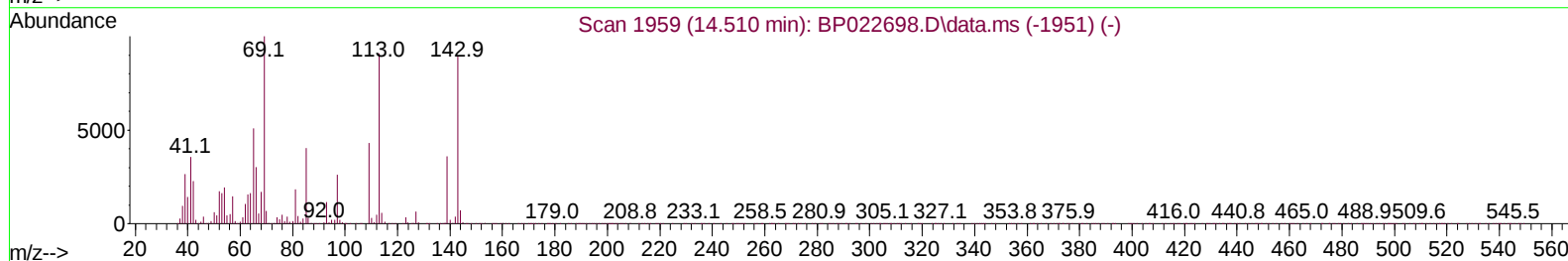
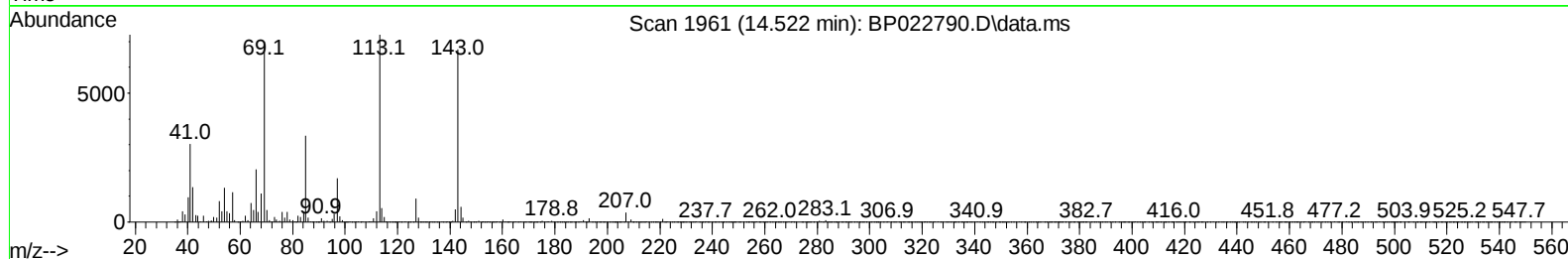
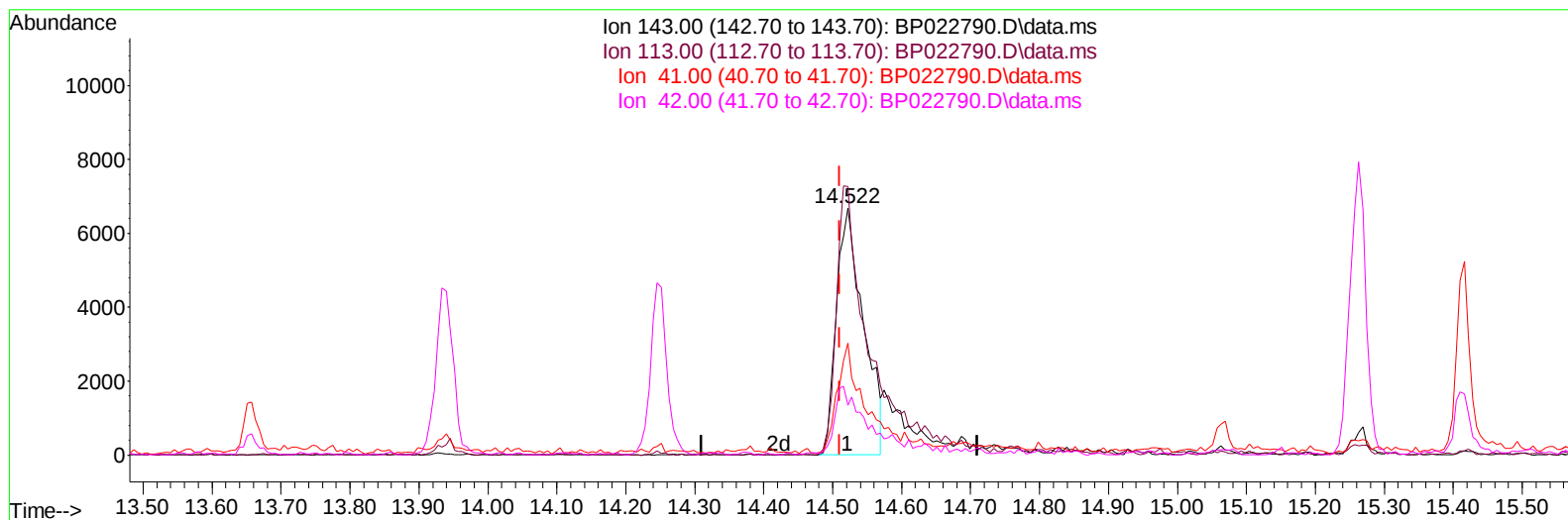
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022790.D
Acq On : 01 Nov 2024 05:13
Operator : RC/JU
Sample : P4630-10
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCP06

Manual IntegrationsAPPROVED

Quant Time: Nov 01 07:56:33 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Oct 30 10:56:39 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/04/2024
Supervised By :mohammad ahmed 11/05/2024



TIC: BP022790.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.012) 5.47 ng/u1

response 18212

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	108.64
41.00	35.90	45.37#
42.00	23.70	20.23

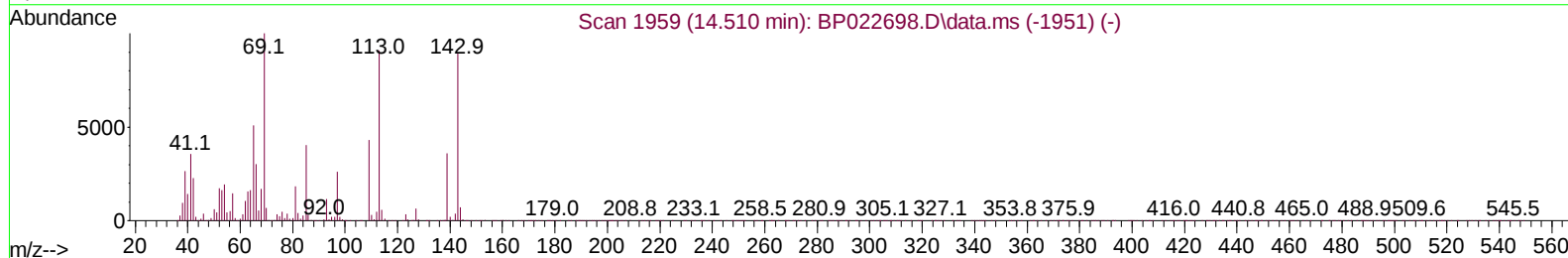
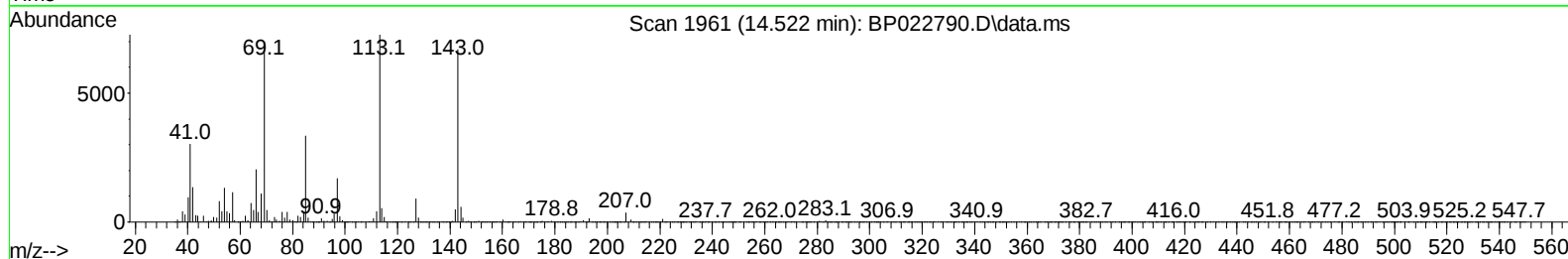
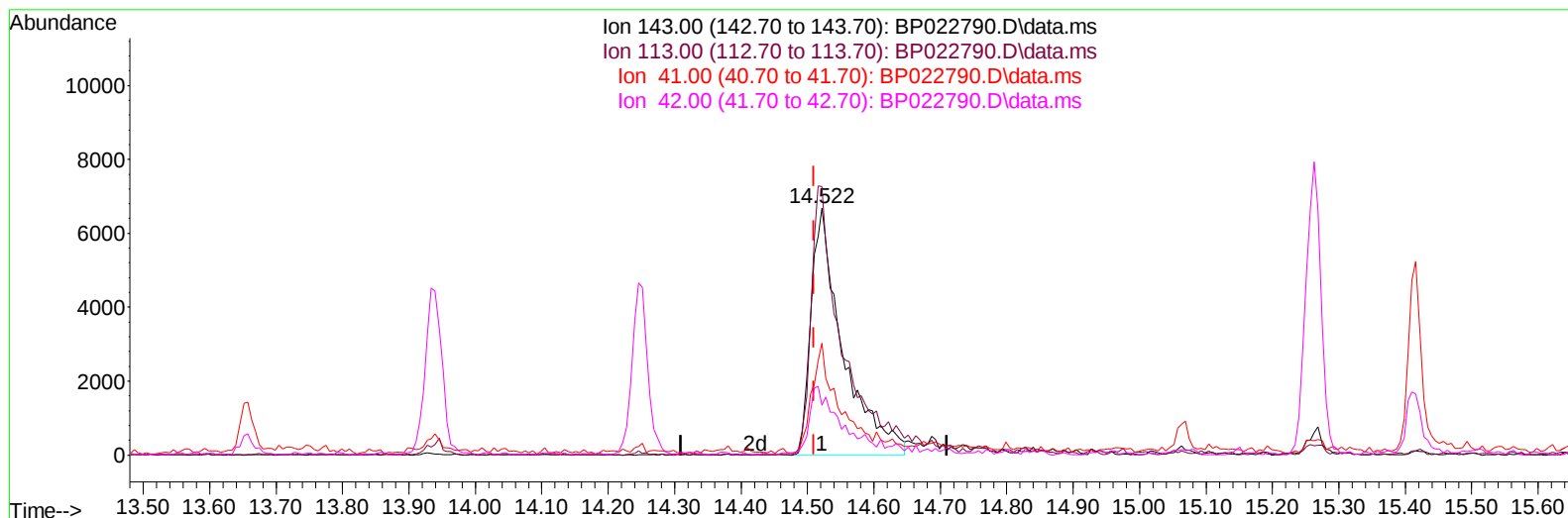
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(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.012) 6.70 ng/ul m

response 22298

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	108.64
41.00	35.90	45.37#
42.00	23.70	20.23

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Quant Time: Nov 01 07:57:18 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.628	152	88072	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.381	136	316982	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.245	164	249618	20.000	ng/ul	-0.02
64) Phenanthrene-d10	17.063	188	623810	20.000	ng/ul	0.00
79) Chrysene-d12	21.510	240	747274	20.000	ng/ul	# 0.01
88) Perylene-d12	24.774	264	904901	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	7782	3.434	ng/uL	0.00
4) Pyridine-d5	3.599	84	38960	6.262	ng/ul	0.00
7) Phenol-d5	6.834	99	57064	7.697	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	115431	26.491	ng/ul	0.00
11) 2-Chlorophenol-d4	7.175	132	135903	25.835	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	108608	18.357	ng/ul	0.00
21) Nitrobenzene-d5	8.775	128	73897	30.320	ng/ul	0.00
24) 2-Nitrophenol-d4	9.487	143	88923	31.514	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.022	165	176750	29.472	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.534	131	142478	20.106	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166	618171	31.169	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	655308	31.109	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.522	143	22298m	6.702	ng/ul	0.01
60) Fluorene-d10	15.263	176	549258	31.929	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.416	200	106709	26.009	ng/ul	0.01
73) Anthracene-d10	17.169	188	998096	34.012	ng/ul	0.00
81) Pyrene-d10	19.557	212	1317765	33.347	ng/ul	0.01
92) Benzo(a)pyrene-d12	24.545	264	1677072	34.703	ng/ul	0.02

Target Compounds Qvalue

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

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