

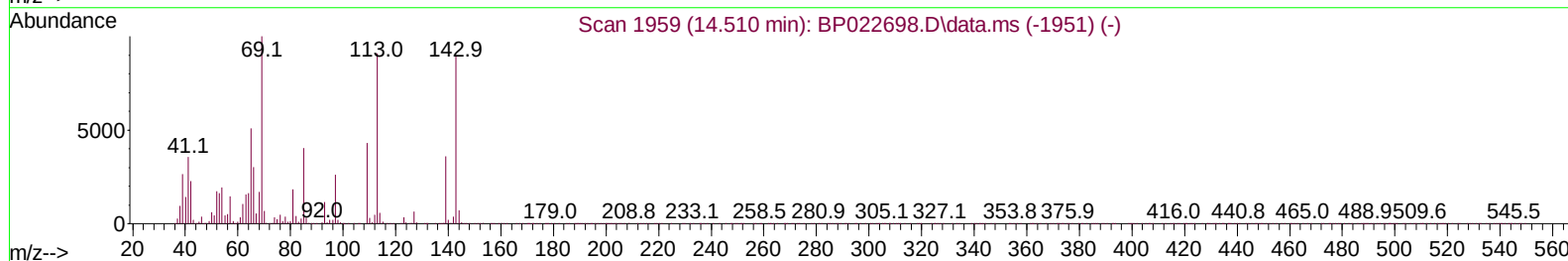
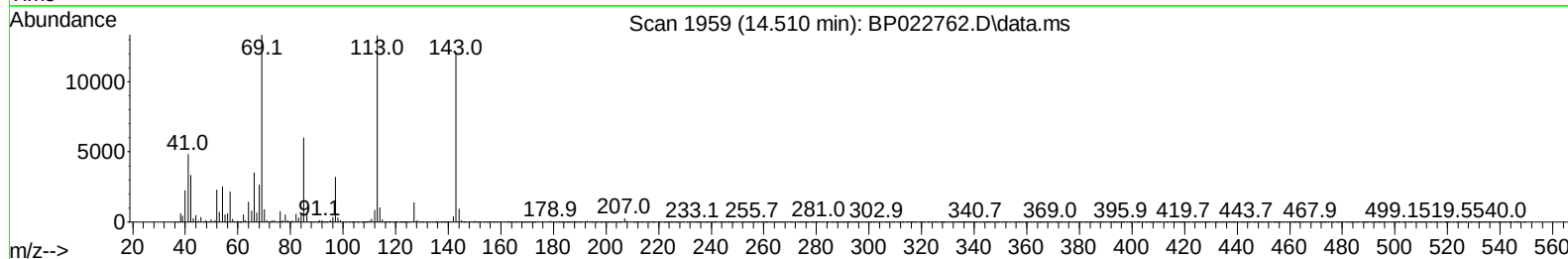
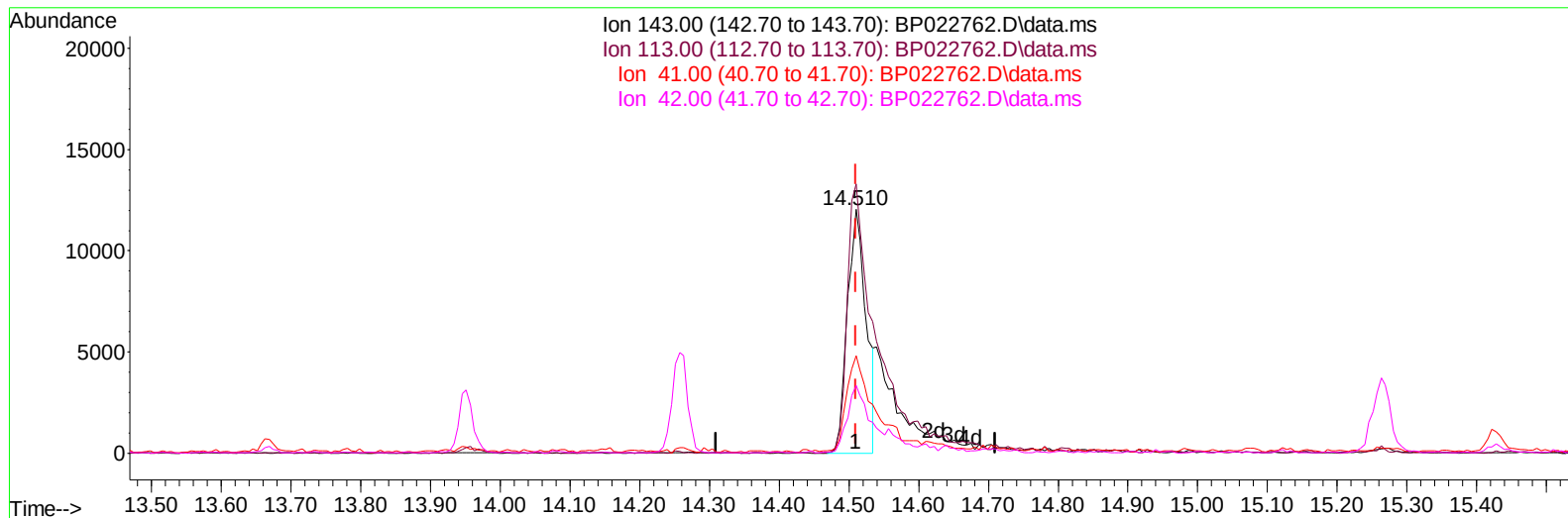
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022762.D
Acq On : 31 Oct 2024 09:22
Operator : RC/JU
Sample : P4592-08
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
F7H08

Manual IntegrationsAPPROVED

Quant Time: Oct 31 12:40:22 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: BP022762.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (-0.000) 6.68 ng/u1

response 22367

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	110.38
41.00	35.90	40.05
42.00	23.70	27.81

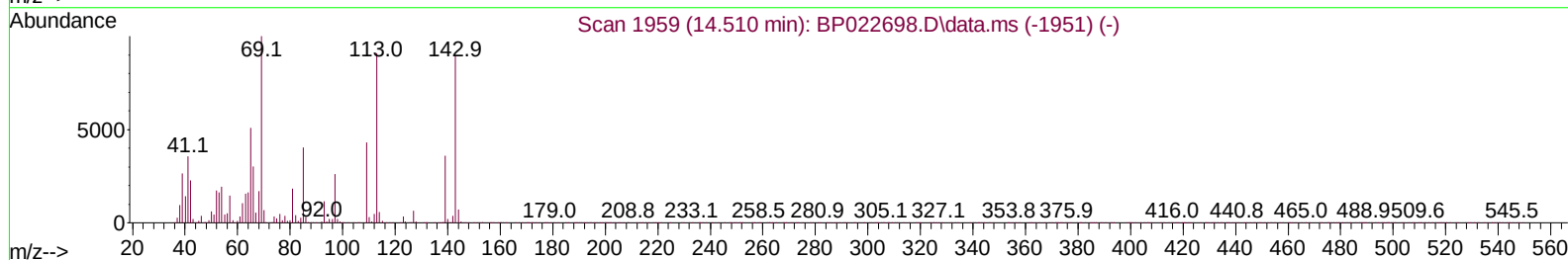
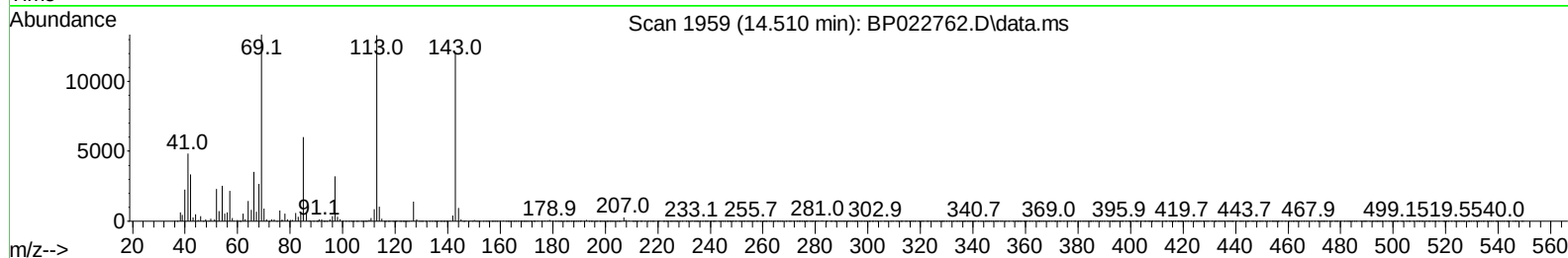
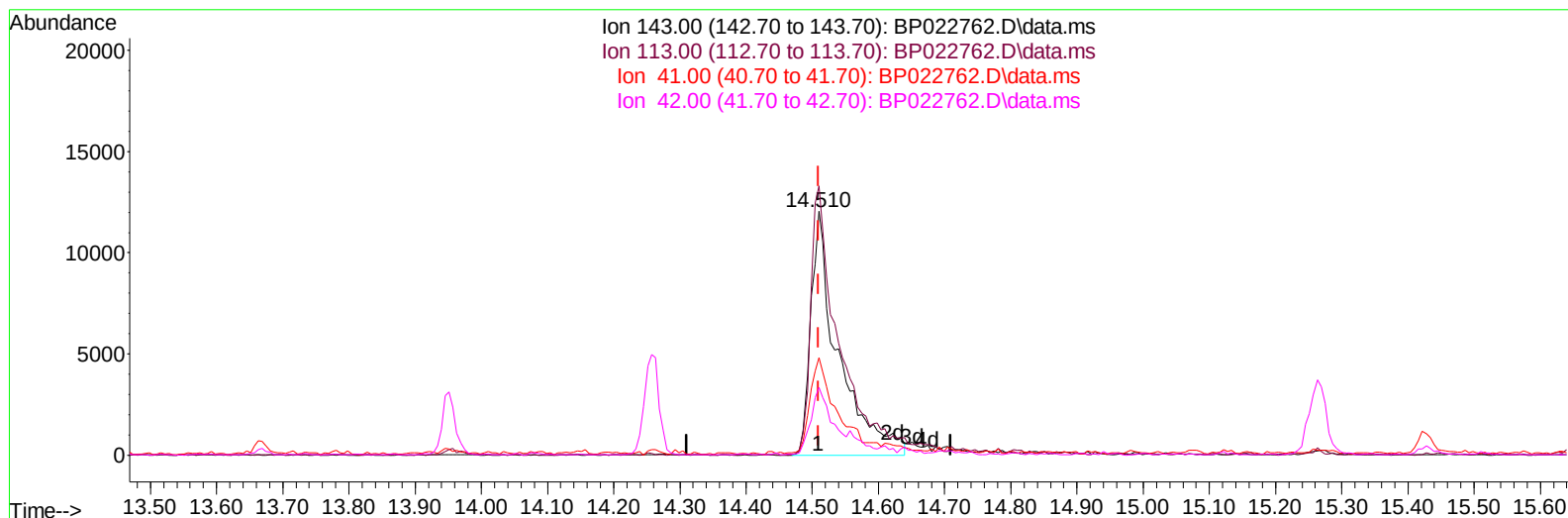
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TIC: BP022762.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (-0.000) 10.46 ng/ul m

response 35013

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	110.38
41.00	35.90	40.05
42.00	23.70	27.81

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Manual IntegrationsAPPROVED

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

Quant Time: Nov 01 01:45:22 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

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	84673	20.000	ng/ul	0.00
20) Naphthalene-d8	10.399	136	322446	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	251100	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.069	188	615418	20.000	ng/ul	0.00
79) Chrysene-d12	21.504	240	696078	20.000	ng/ul	# 0.00
88) Perylene-d12	24.762	264	824369	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	5517	2.532	ng/uL	0.00
4) Pyridine-d5	3.593	84	72567	12.132	ng/ul	0.00
7) Phenol-d5	6.828	99	103231	14.483	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	59090	14.105	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	77247	15.274	ng/ul	0.00
15) 4-Methylphenol-d8	8.346	113	85484	15.029	ng/ul	0.00
21) Nitrobenzene-d5	8.775	128	37035	14.938	ng/ul	0.00
24) 2-Nitrophenol-d4	9.487	143	46015	16.031	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.040	165	95386	15.635	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	105034	14.571	ng/ul	0.00
46) Dimethylphthalate-d6	13.669	166	337422	16.913	ng/ul	0.00
49) Acenaphthylene-d8	13.951	160	347831	16.415	ng/ul	0.00
54) 4-Nitrophenol-d4	14.510	143	35013m	10.461	ng/ul	0.00
60) Fluorene-d10	15.263	176	299337	17.298	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.428	200	29315	7.243	ng/ul	0.02
73) Anthracene-d10	17.175	188	518172	17.898	ng/ul	0.00
81) Pyrene-d10	19.545	212	644813	17.517	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.527	264	777061	17.650	ng/ul	0.00

Target Compounds Qvalue

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

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