

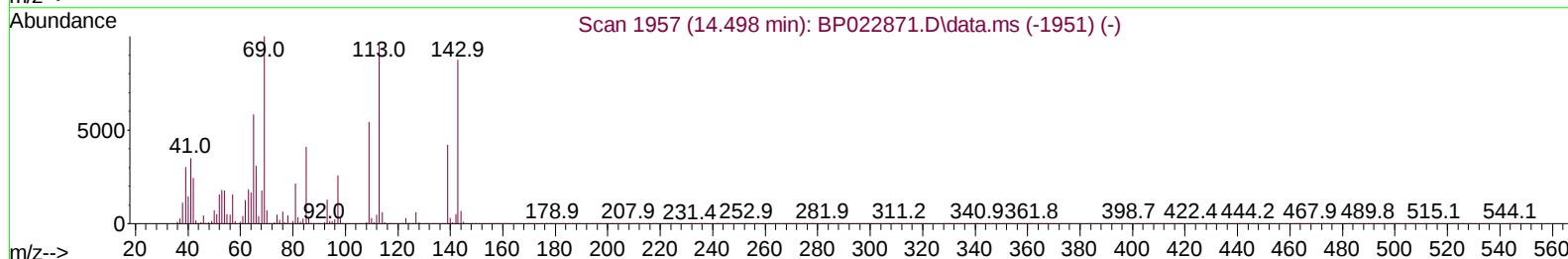
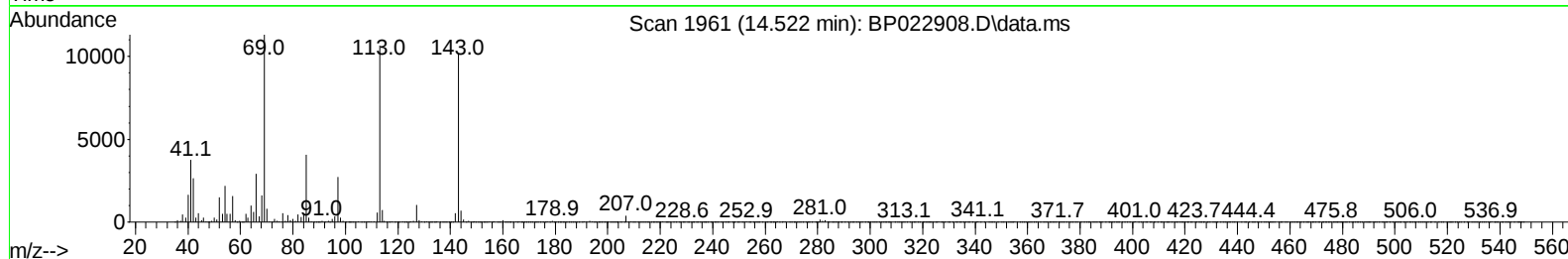
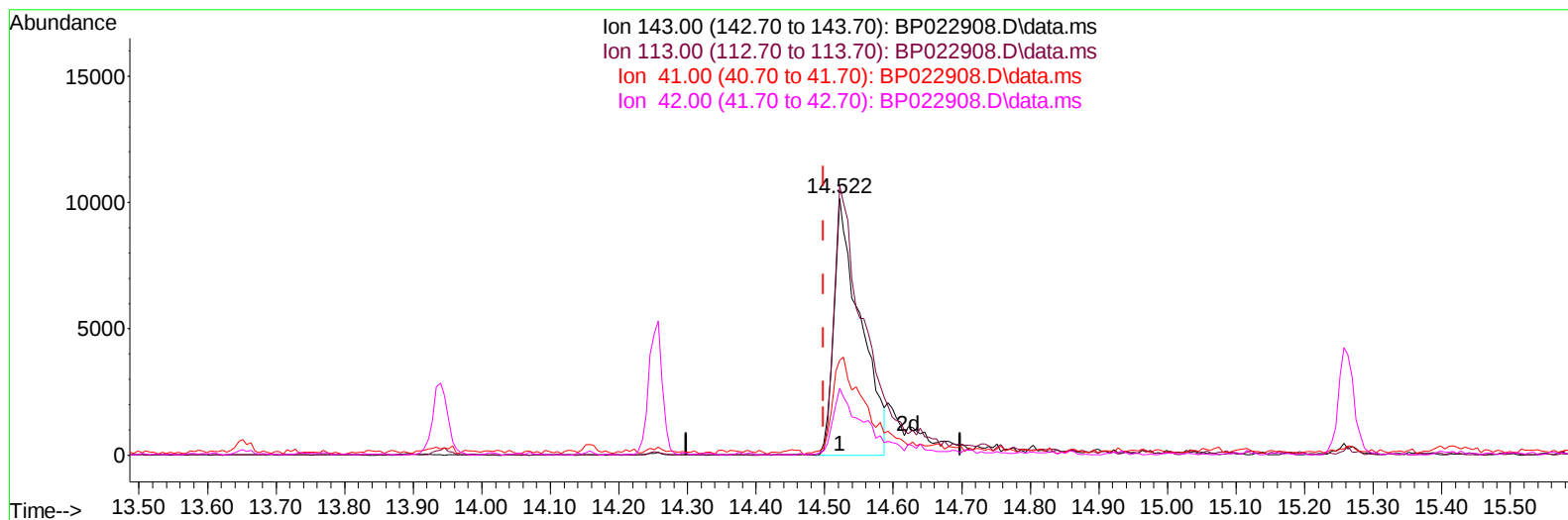
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
 Data File : BP022908.D
 Acq On : 07 Nov 2024 00:43
 Operator : RC/JU
 Sample : P4622-09
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH6N5

Manual IntegrationsAPPROVED

Quant Time: Nov 07 01:20:54 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Nov 06 00:31:03 2024
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/07/2024
 Supervised By :mohammad ahmed 11/08/2024



TIC: BP022908.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.024) 7.61 ng/u1

response 27122

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	104.78
41.00	35.90	36.92
42.00	23.70	26.01

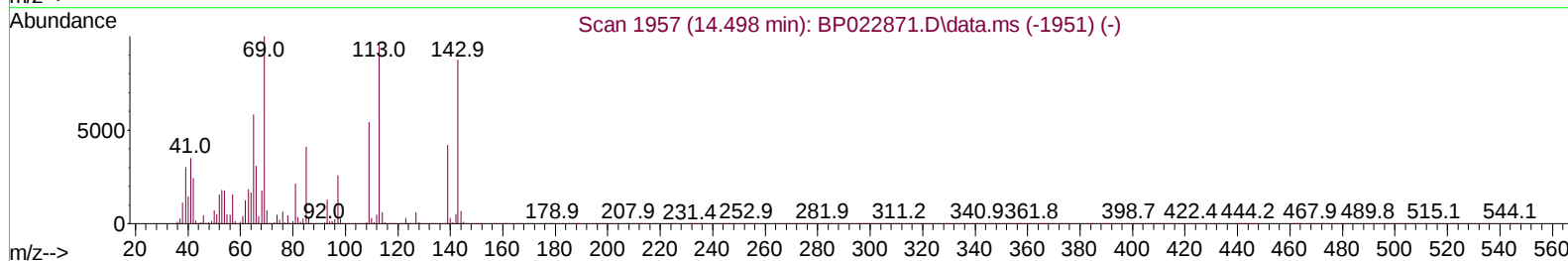
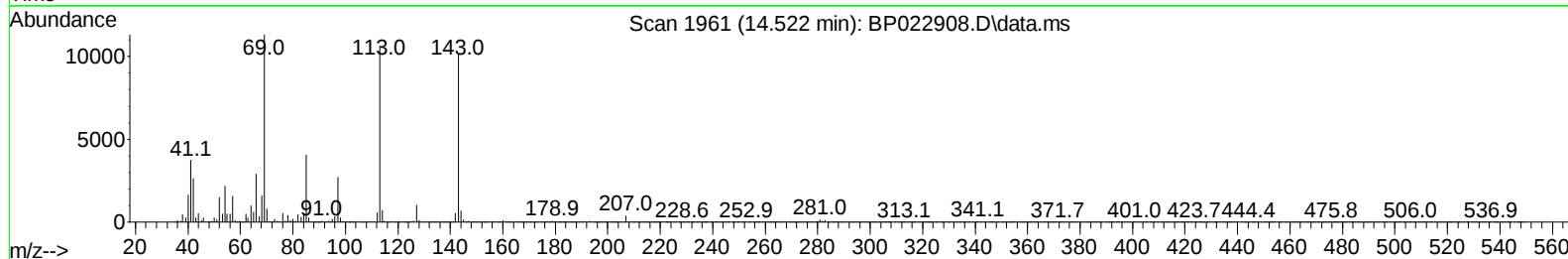
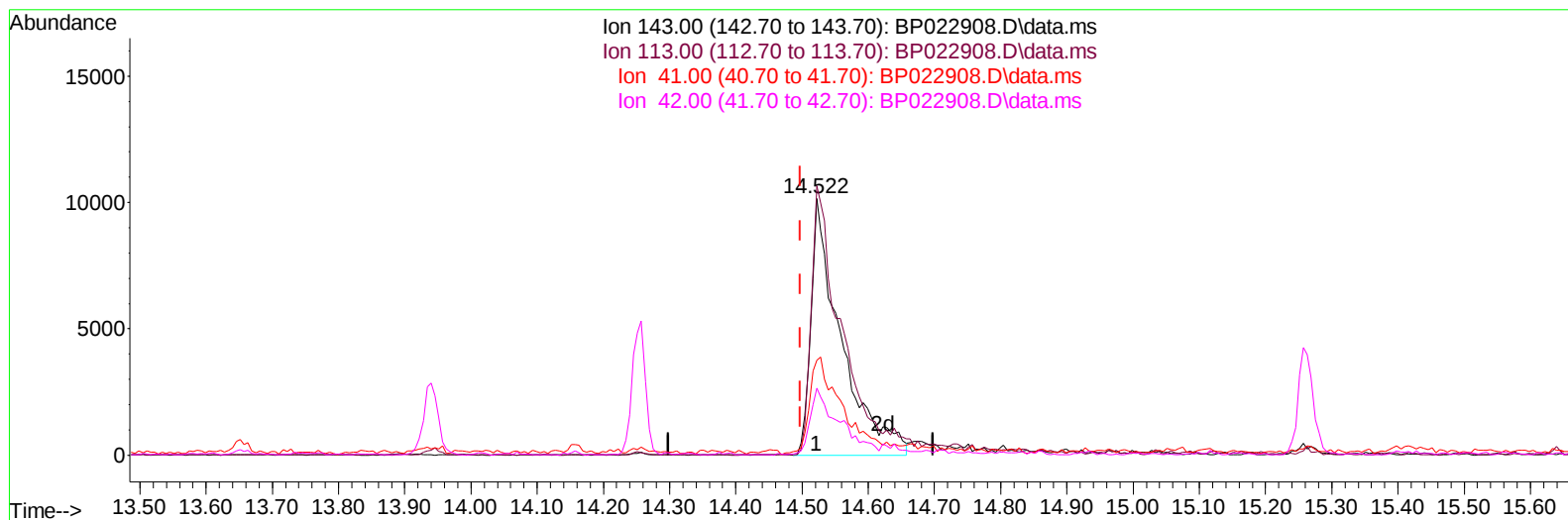
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TIC: BP022908.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.024) 8.93 ng/ul m

response 31818

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	104.78
41.00	35.90	36.92
42.00	23.70	26.01

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 11/07/2024
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Quant Time: Nov 07 01:22:11 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	87715	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.381	136	338708	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.257	164	267451	20.000	ng/ul	0.02
64) Phenanthrene-d10	17.057	188	647995	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	682099	20.000	ng/ul	0.00
88) Perylene-d12	24.745	264	834741	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	4061	1.799	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	6.822	99	102855	13.930	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.964	67	58964	13.587	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	77750	14.840	ng/ul	0.00
15) 4-Methylphenol-d8	8.328	113	83408	14.155	ng/ul	0.00
21) Nitrobenzene-d5	8.769	128	40036	15.373	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	43230	14.338	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.028	165	98778	15.414	ng/ul	0.01
31) 4-Chloroaniline-d4	10.528	131	88092	11.634	ng/ul	0.00
46) Dimethylphthalate-d6	13.651	166	250111	11.770	ng/ul	-0.01
49) Acenaphthylene-d8	13.940	160	376280	16.672	ng/ul	0.00
54) 4-Nitrophenol-d4	14.522	143	31818m	8.925	ng/ul	0.02
60) Fluorene-d10	15.263	176	315308	17.107	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.416	200	11081	2.600	ng/ul	0.00
73) Anthracene-d10	17.151	188	518572	17.012	ng/ul	-0.01
81) Pyrene-d10	19.545	212	657346	18.224	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.521	264	765165	17.164	ng/ul	0.00

Target Compounds Qvalue

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

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