

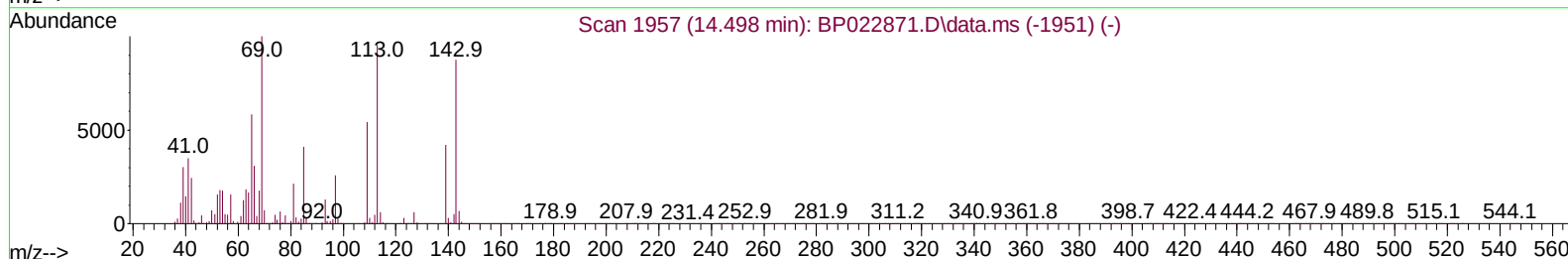
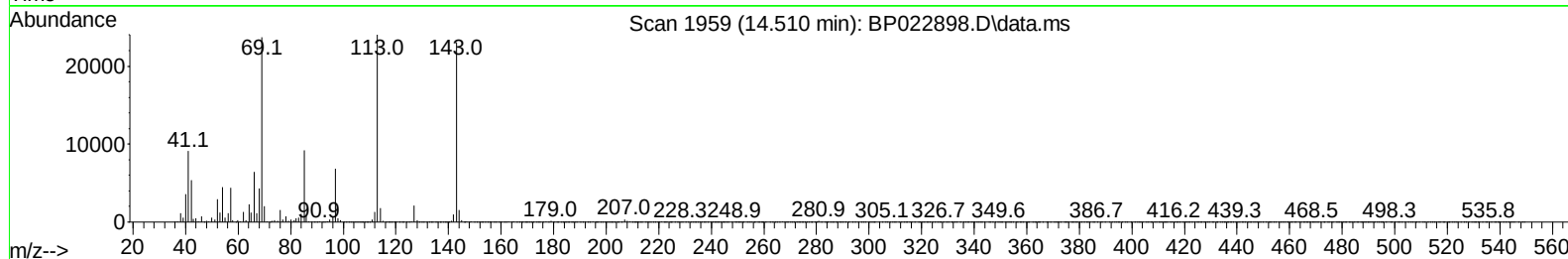
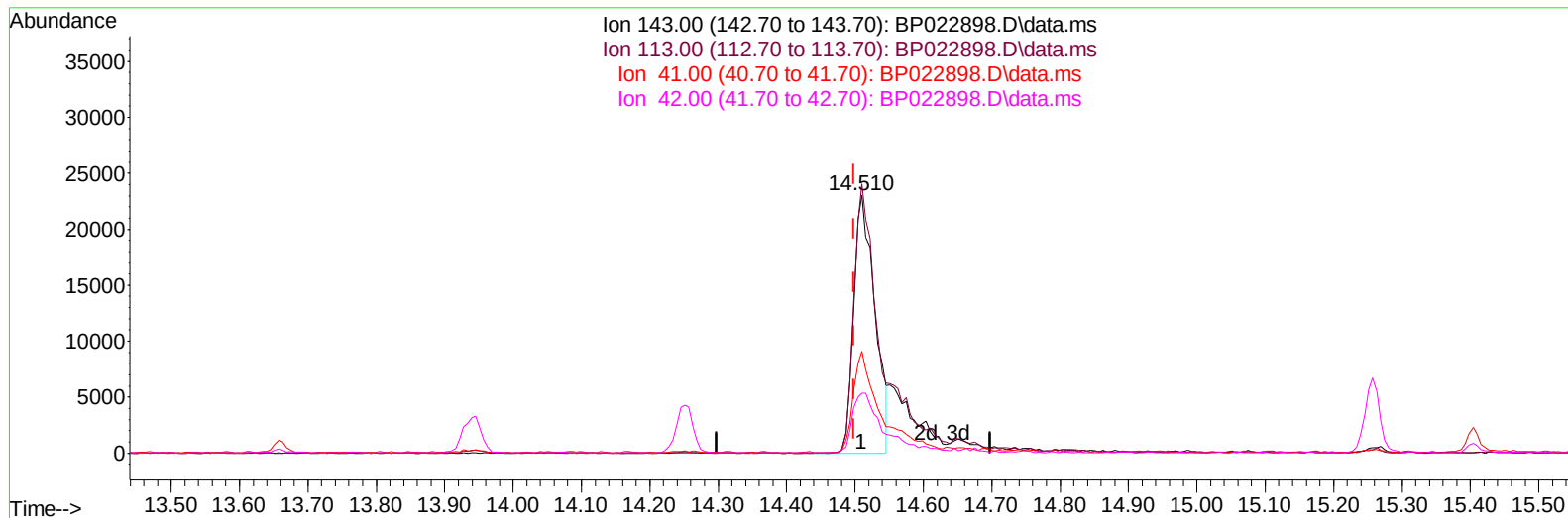
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022898.D
Acq On : 06 Nov 2024 17:55
Operator : RC/JU
Sample : PB164642BL
Misc :
ALS Vial : 45 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
SBLK642

Manual IntegrationsAPPROVED

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

Quant Time: Nov 07 00:42:42 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



TIC: BP022898.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.012) 14.08 ng/u1

response 49574

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	104.31
41.00	35.90	39.52
42.00	23.70	23.20

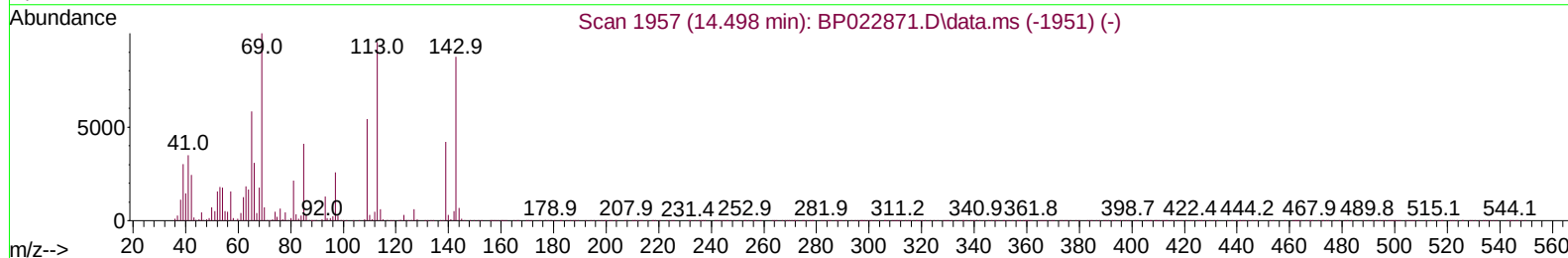
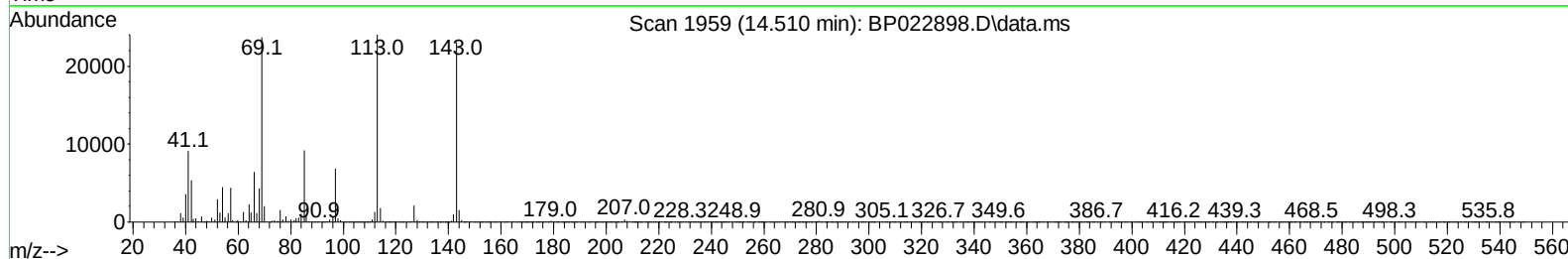
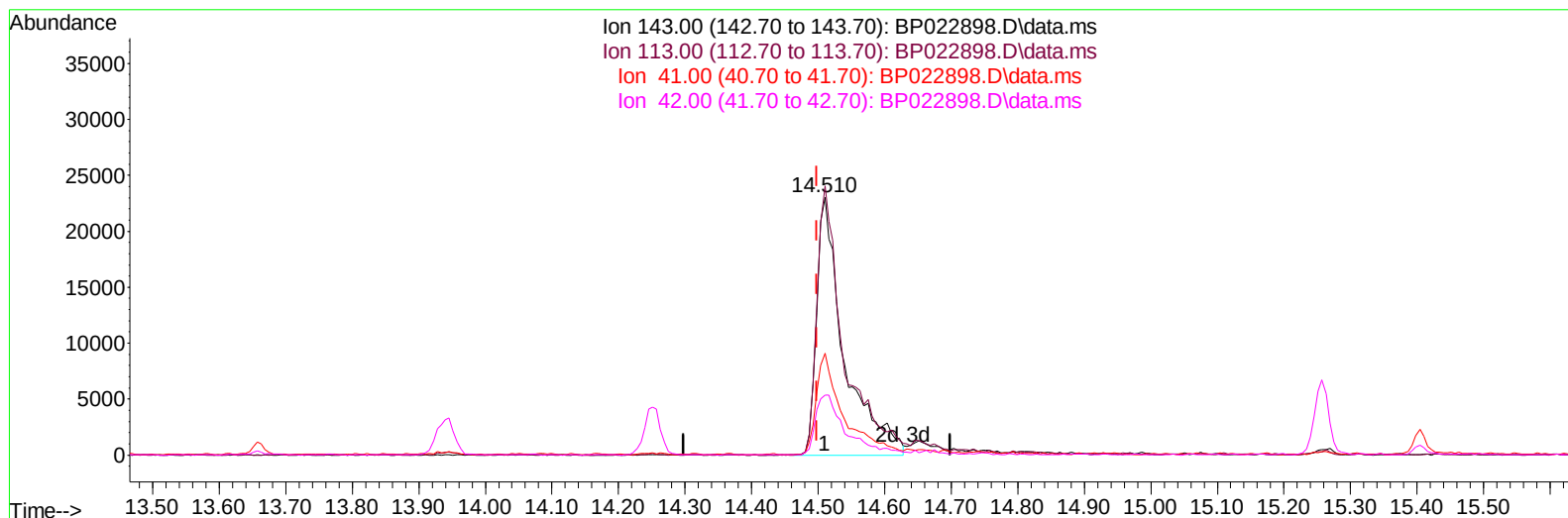
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TIC: BP022898.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.012) 18.67 ng/ul m

response 65709

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	104.31
41.00	35.90	39.52
42.00	23.70	23.20

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Reviewed By :Jagrut Upadhyay 11/07/2024
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Quant Time: Nov 07 00:45:23 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.616	152	86717	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.375	136	333898	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	264088	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.057	188	625670	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240	641804	20.000	ng/ul	-0.01
88) Perylene-d12	24.733	264	735235	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.170	96	8351	3.742	ng/uL	0.00
4) Pyridine-d5	3.587	84	120400	19.654	ng/ul	0.00
7) Phenol-d5	6.822	99	166400	22.796	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.963	67	90326	21.054	ng/ul	0.00
11) 2-Chlorophenol-d4	7.163	132	126538	24.430	ng/ul	0.00
15) 4-Methylphenol-d8	8.328	113	138132	23.712	ng/ul	0.00
21) Nitrobenzene-d5	8.769	128	62910	24.504	ng/ul	0.00
24) 2-Nitrophenol-d4	9.475	143	76233	25.648	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.016	165	153167	24.246	ng/ul	0.00
31) 4-Chloroaniline-d4	10.522	131	170306	22.815	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166	508854	24.251	ng/ul	0.00
49) Acenaphthylene-d8	13.940	160	552092	24.773	ng/ul	0.00
54) 4-Nitrophenol-d4	14.510	143	65709m	18.667	ng/ul	0.01
60) Fluorene-d10	15.257	176	450481	24.752	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	50996	12.393	ng/ul	-0.01
73) Anthracene-d10	17.151	188	770234	26.169	ng/ul	-0.01
81) Pyrene-d10	19.533	212	951386	28.032	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.515	264	1044865	26.610	ng/ul	-0.01

Target Compounds Qvalue

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

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