

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022904.D
Acq On : 06 Nov 2024 22:01
Operator : RC/JU
Sample : P4622-05
Misc :
ALS Vial : 51 Sample Multiplier: 1

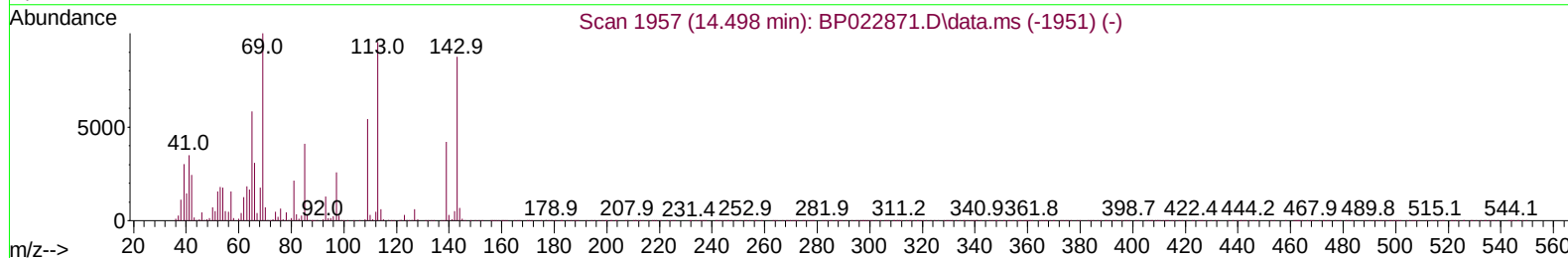
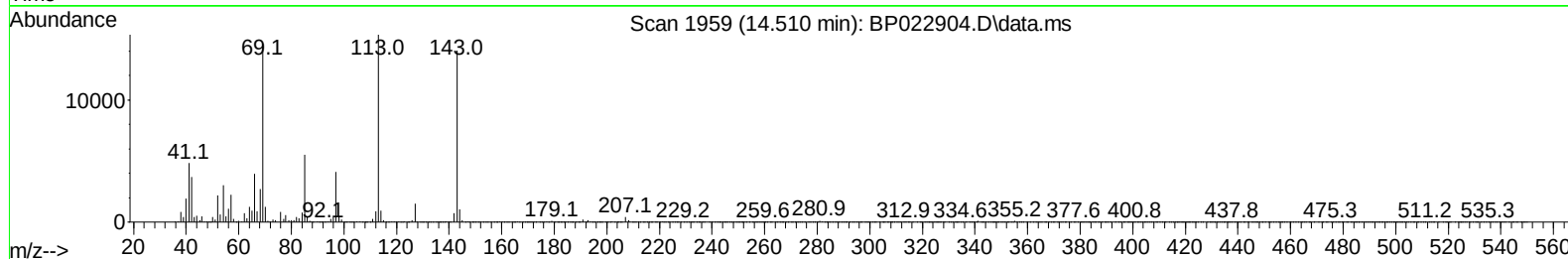
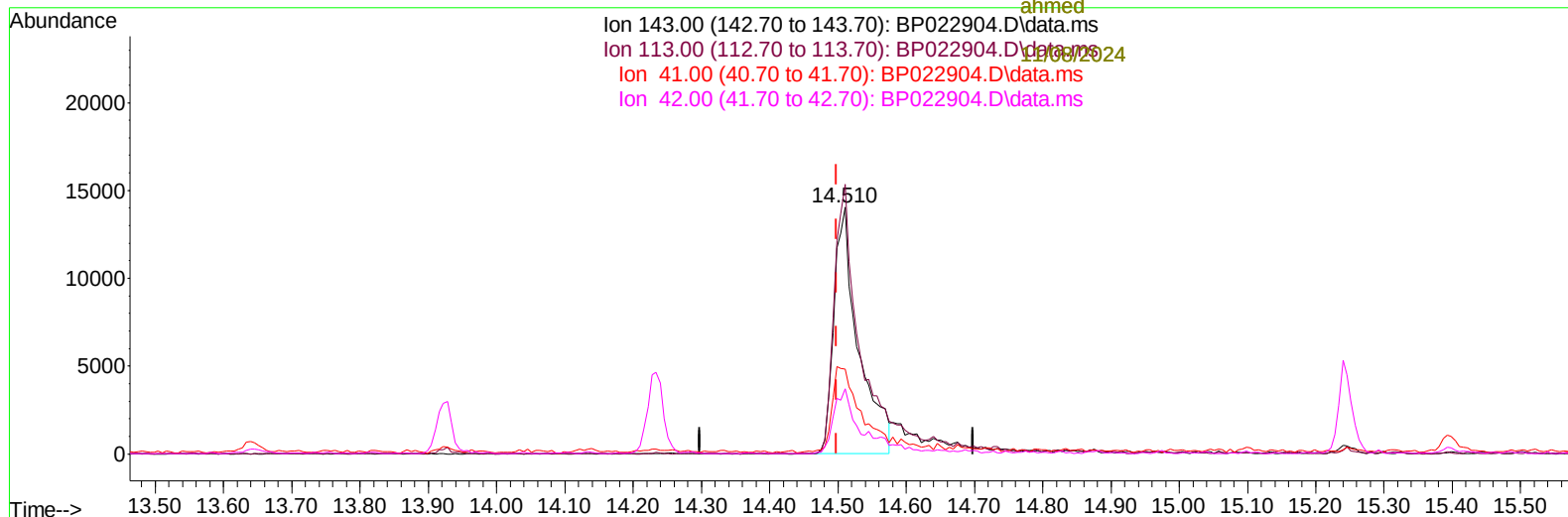
Instrument :
BNA_P
ClientSampleId :
BH6N1

Manual IntegrationsAPPROVED

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

Quant Time: Nov 07 01:09:33 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

11/07/2024
Supervised By :mohammad
ahmed



TIC: BP022904.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.012) 9.92 ng/ul

response 35135

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	109.20
41.00	35.90	34.27
42.00	23.70	26.31

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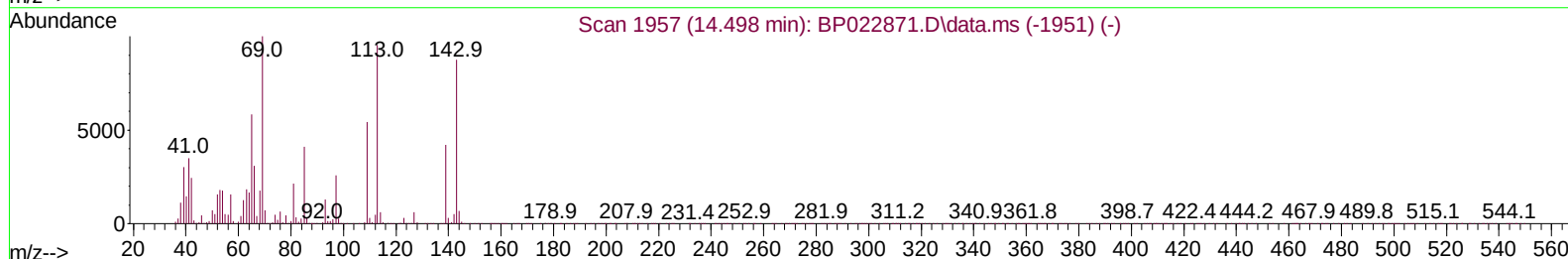
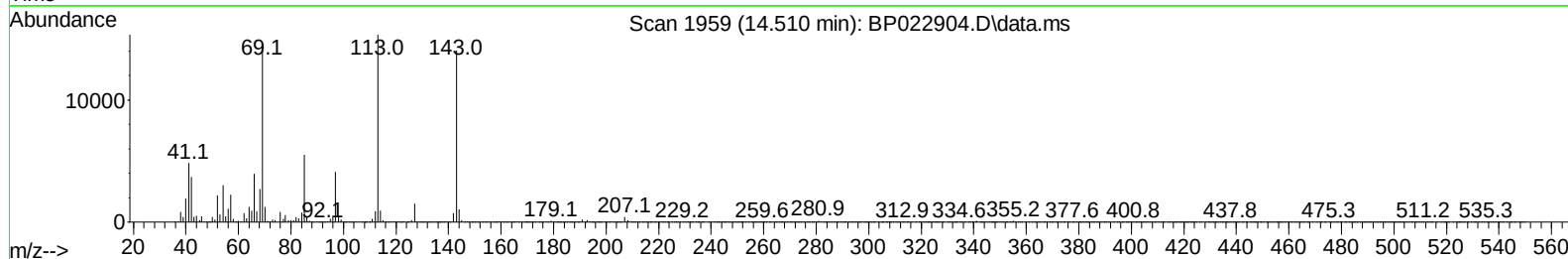
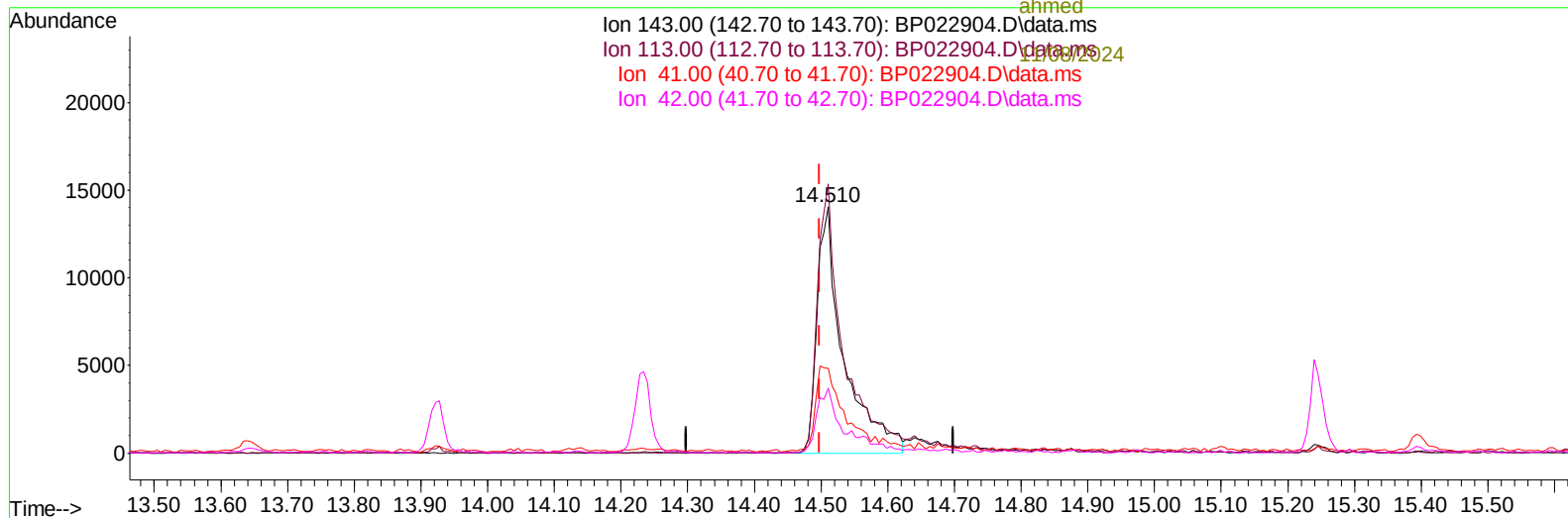
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TIC: BP022904.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.012) 10.96 ng/ul m

response 38822

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	109.20
41.00	35.90	34.27
42.00	23.70	26.31

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----11/08/2024						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.622	152	89660	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	342561	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.234	164	265851	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.040	188	633480	20.000	ng/ul	-0.01
79) Chrysene-d12	21.492	240	674785	20.000	ng/ul	# 0.00
88) Perylene-d12	24.751	264	805866	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.176	96	4546	1.970	ng/uL	0.00
4) Pyridine-d5	3.593	84	36475	5.759	ng/ul	0.00
7) Phenol-d5	6.822	99	107200	14.204	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.964	67	59937	13.512	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	81885	15.290	ng/ul	0.00
15) 4-Methylphenol-d8	8.328	113	91837	15.247	ng/ul	0.00
21) Nitrobenzene-d5	8.763	128	41236	15.656	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	48930	16.046	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	102157	15.762	ng/ul	0.00
31) 4-Chloroaniline-d4	10.522	131	115968	15.143	ng/ul	0.00
46) Dimethylphthalate-d6	13.640	166	354816	16.798	ng/ul	-0.02
49) Acenaphthylene-d8	13.922	160	393568	17.543	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.510	143	38822m	10.956	ng/ul	0.01
60) Fluorene-d10	15.245	176	325676	17.776	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.398	200	21826	5.239	ng/ul	-0.02
73) Anthracene-d10	17.139	188	512287	17.191	ng/ul	-0.02
81) Pyrene-d10	19.522	212	673809	18.883	ng/ul	-0.02
92) Benzo(a)pyrene-d12	24.527	264	785160	18.244	ng/ul	0.00

Target Compounds Qvalue

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

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