

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022902.D
Acq On : 06 Nov 2024 20:39
Operator : RC/JU
Sample : P4622-03
Misc :
ALS Vial : 49 Sample Multiplier: 1

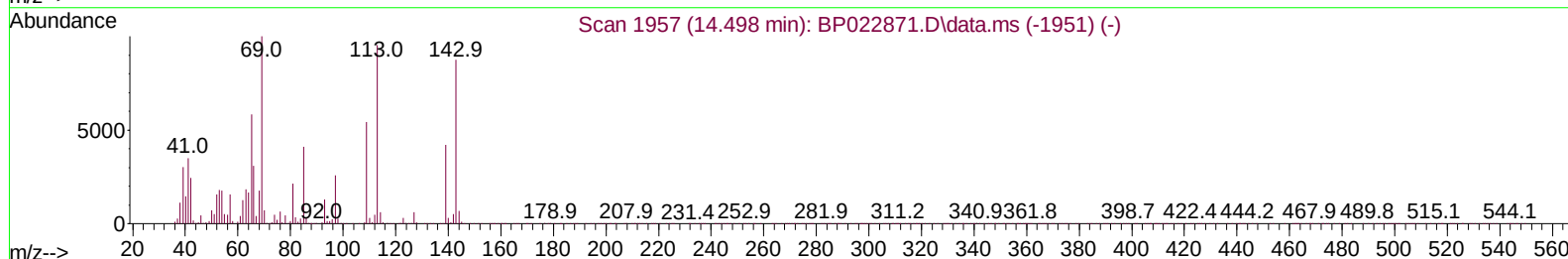
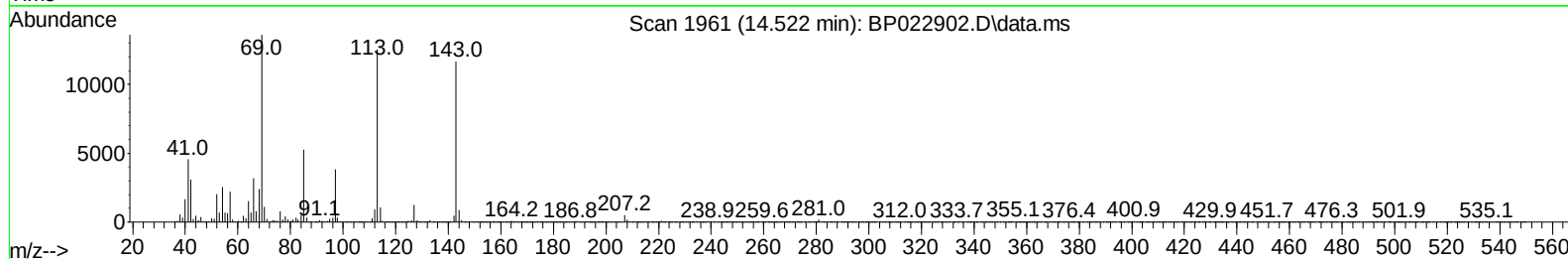
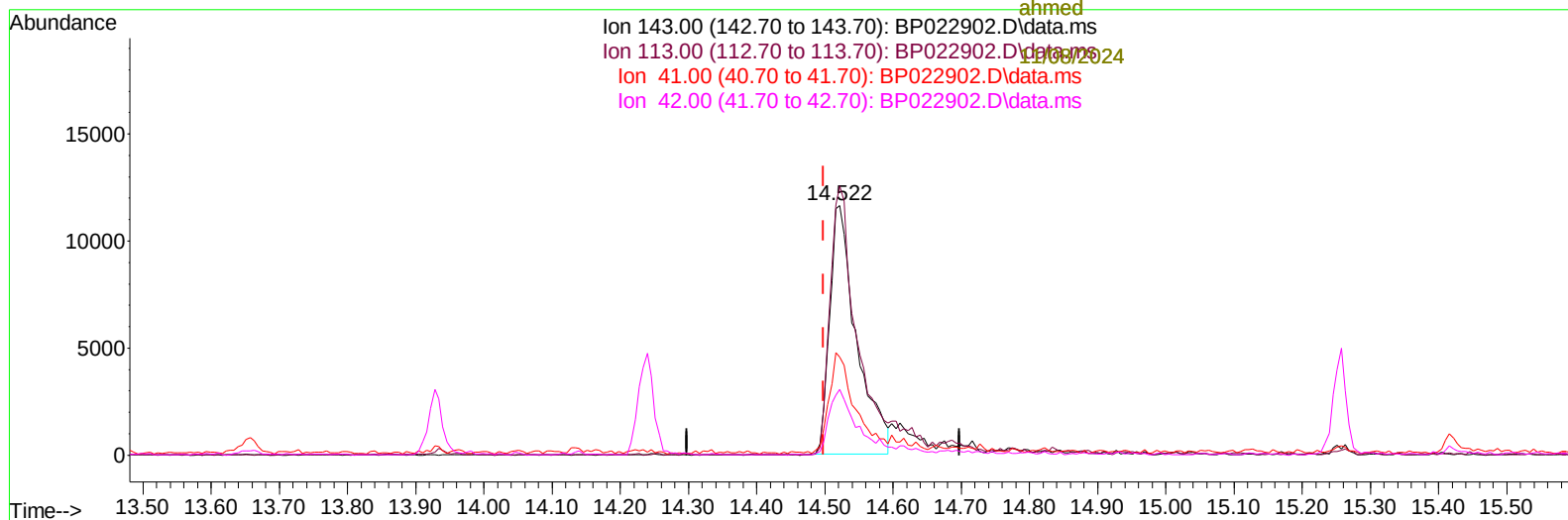
Instrument :
BNA_P
ClientSampleId :
BH6M9

Manual IntegrationsAPPROVED

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

Quant Time: Nov 07 01:03:08 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: BP022902.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.024) 9.52 ng/u1

response 31741

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	107.76
41.00	35.90	39.16
42.00	23.70	26.40

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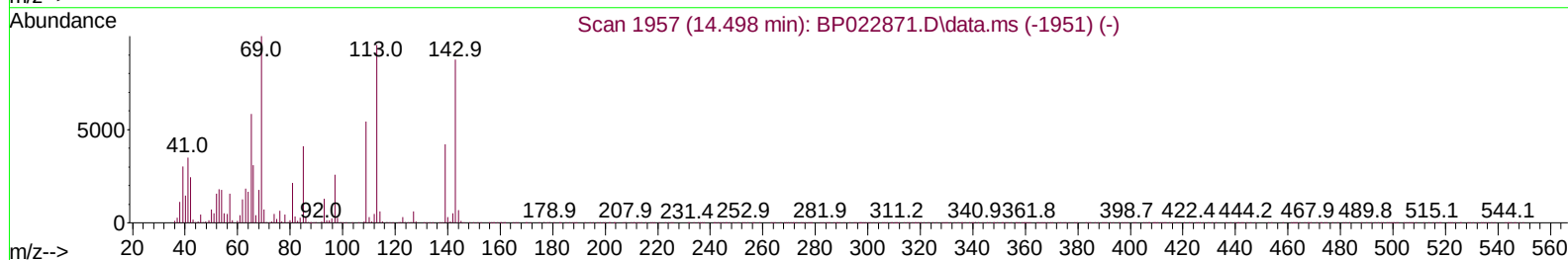
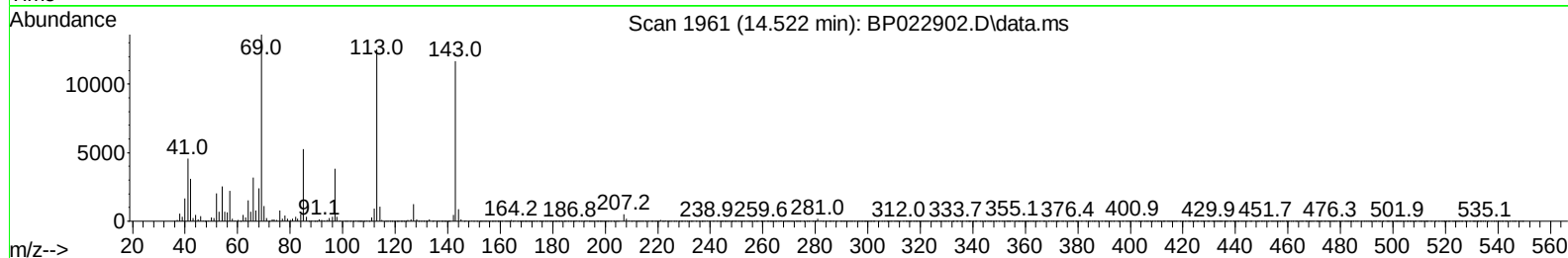
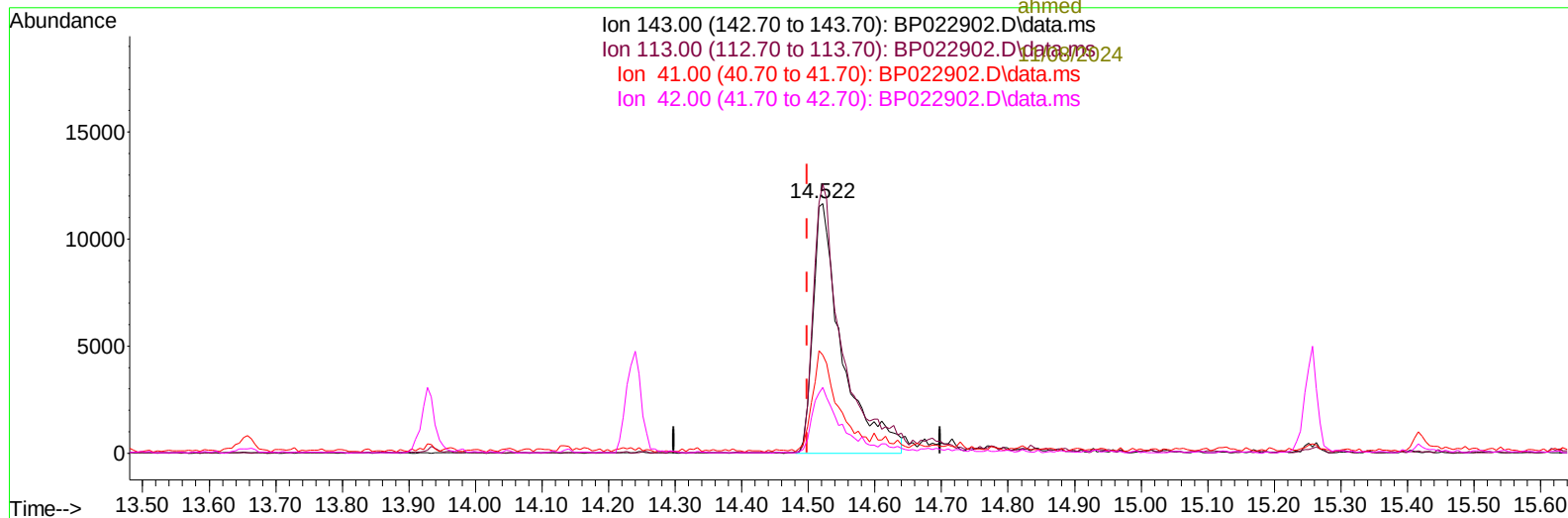
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ahmed



TIC: BP022902.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.522min (+ 0.024) 10.56 ng/ul m

response 35216

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	107.76
41.00	35.90	39.16
42.00	23.70	26.40

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Reviewed By :Jagrut Upadhyay 11/07/2024
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Quant Time: Nov 07 01:05:33 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

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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	83373	20.000	ng/ul	0.00
20) Naphthalene-d8	10.375	136	325596	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.240	164	250161	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.045	188	592200	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	626130	20.000	ng/ul	0.00
88) Perylene-d12	24.739	264	750814	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	4074	1.899	ng/uL	0.00
4) Pyridine-d5	3.593	84	52156	8.855	ng/ul	0.00
7) Phenol-d5	6.816	99	91112	12.982	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.963	67	49700	12.049	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	68362	13.728	ng/ul	0.00
15) 4-Methylphenol-d8	8.334	113	74855	13.365	ng/ul	0.00
21) Nitrobenzene-d5	8.769	128	34467	13.767	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	42975	14.827	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	90017	14.613	ng/ul	0.00
31) 4-Chloroaniline-d4	10.528	131	99204	13.629	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166	316498	15.924	ng/ul	0.00
49) Acenaphthylene-d8	13.928	160	336615	15.945	ng/ul	0.00
54) 4-Nitrophenol-d4	14.522	143	35216m	10.561	ng/ul	0.02
60) Fluorene-d10	15.257	176	285054	16.535	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.416	200	22273	5.719	ng/ul	0.00
73) Anthracene-d10	17.151	188	453772	16.288	ng/ul	-0.01
81) Pyrene-d10	19.528	212	587014	17.729	ng/ul	-0.01
92) Benzo(a)pyrene-d12	24.510	264	681694	17.001	ng/ul	-0.02

Target Compounds Qvalue

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

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