

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP103024\
Data File : BP022756.D
Acq On : 31 Oct 2024 05:19
Operator : RC/JU
Sample : P4592-02
Misc :
ALS Vial : 23 Sample Multiplier: 1

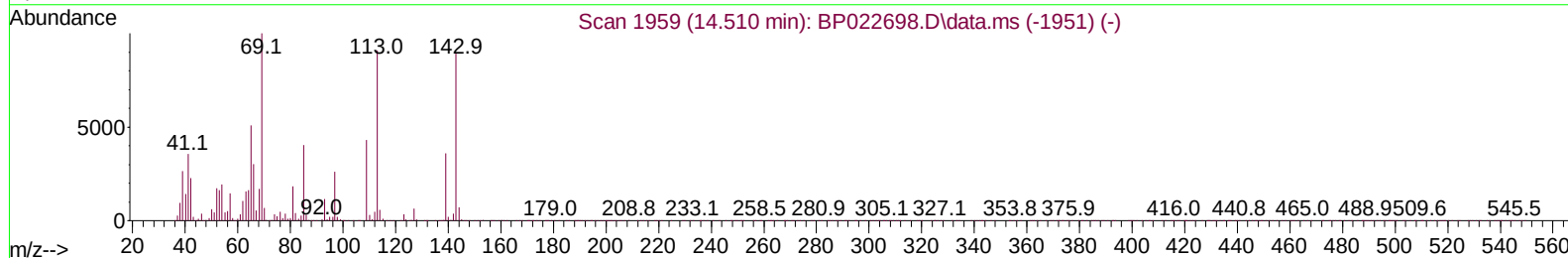
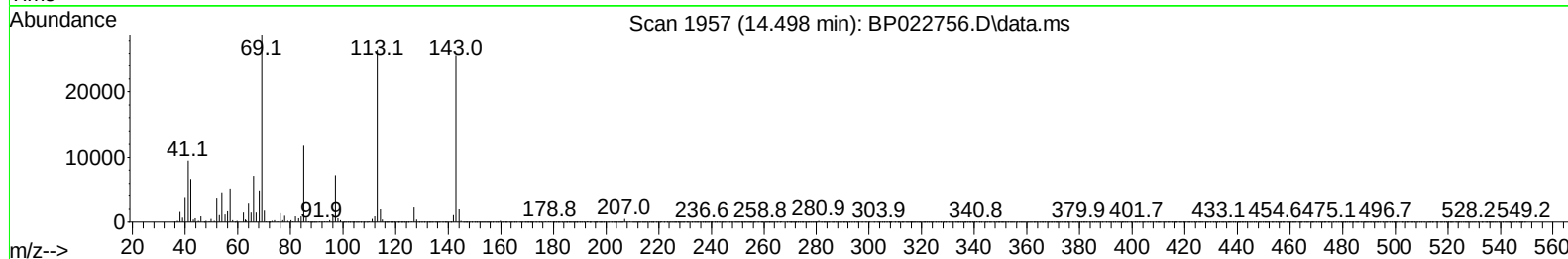
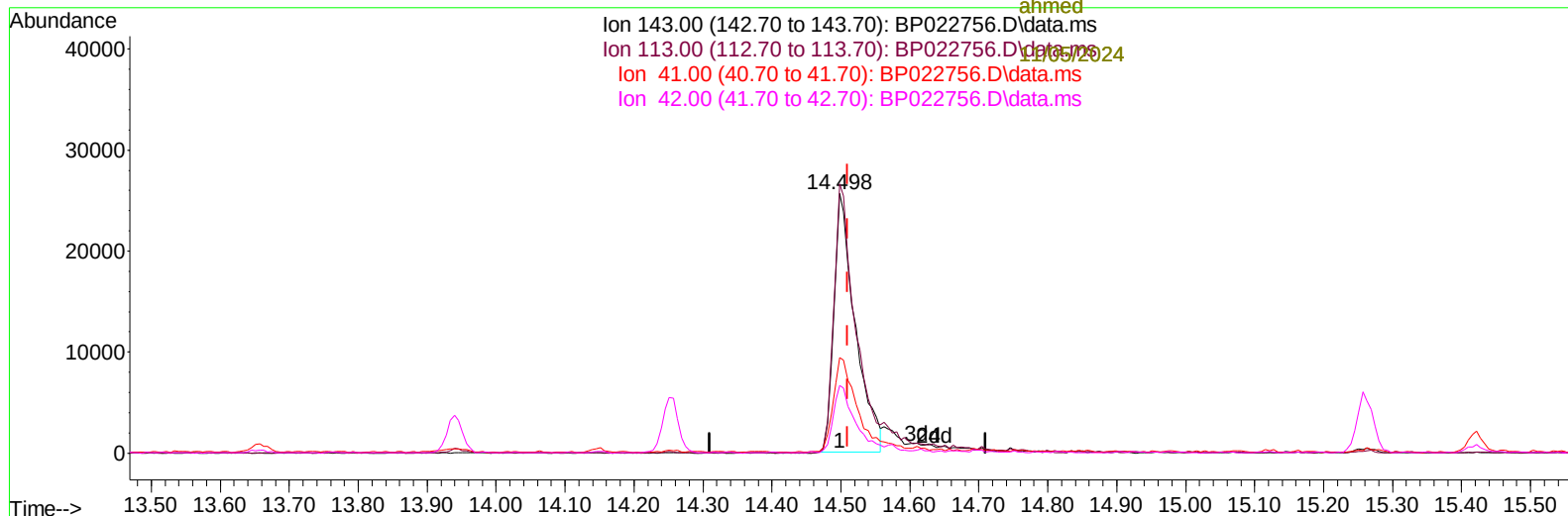
Instrument :
BNA_P
ClientSampleId :
F7H02

Manual IntegrationsAPPROVED

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

11/04/2024
Supervised By :mohammad
ahmed

Quant Time: Oct 31 12:39:19 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



TIC: BP022756.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.012) 13.98 ng/u1

response 55123

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	103.67
41.00	35.90	36.87
42.00	23.70	26.02

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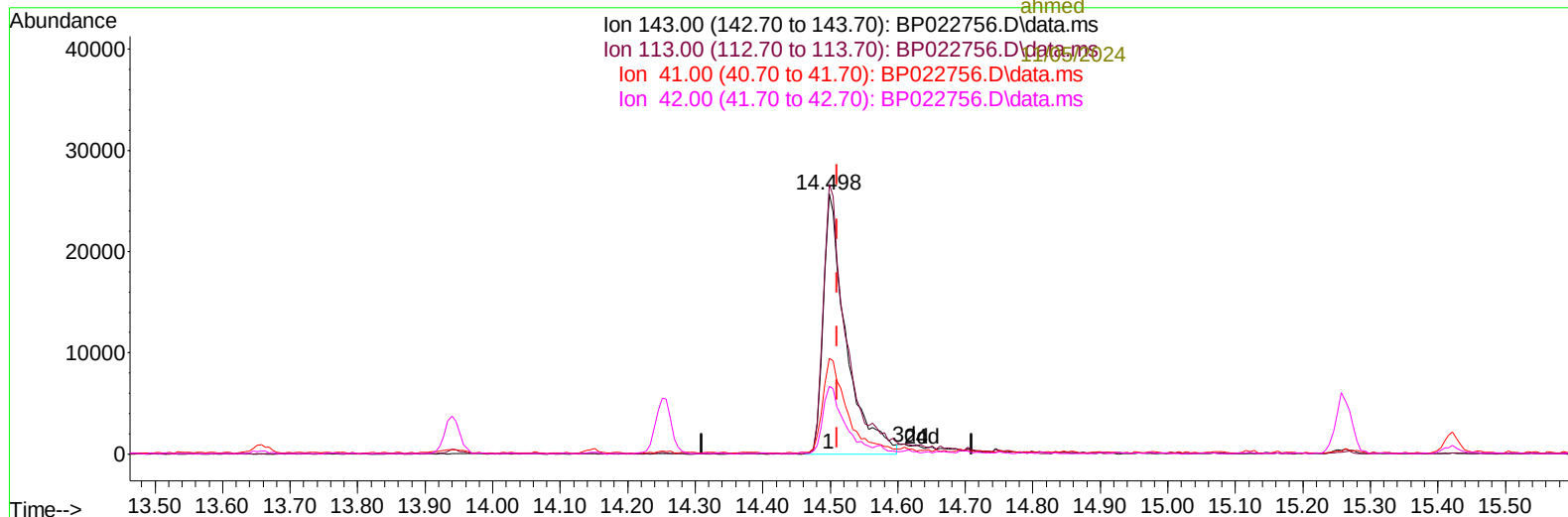
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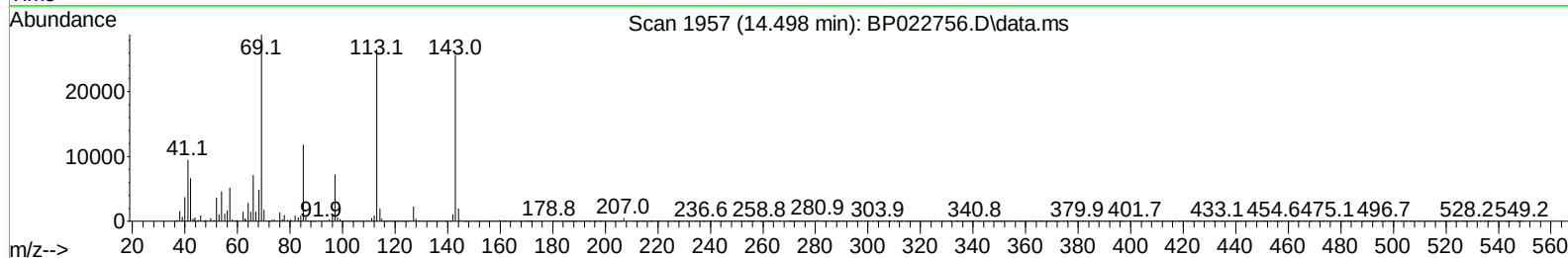
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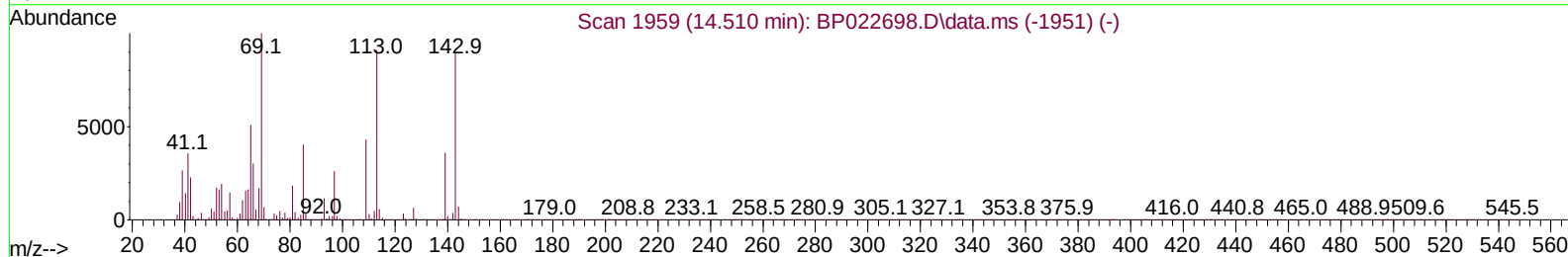
Ion 143.00 (142.70 to 143.70): BP022756.D\data.ms
Ion 113.00 (112.70 to 113.70): BP022756.D\data.ms
Ion 41.00 (40.70 to 41.70): BP022756.D\data.ms
Ion 42.00 (41.70 to 42.70): BP022756.D\data.ms



Scan 1957 (14.498 min): BP022756.D\data.ms



Scan 1959 (14.510 min): BP022698.D\data.ms (-1951) (-)



TIC: BP022756.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.012) 15.17 ng/ul m

response 59799

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	97.60	103.67
41.00	35.90	36.87
42.00	23.70	26.02

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Quant Time: Nov 01 01:17:08 2024
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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----11/05/2024						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	95703	20.000	ng/ul	0.00
20) Naphthalene-d8	10.387	136	369506	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.251	164	295775	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.057	188	730185	20.000	ng/ul	0.00
79) Chrysene-d12	21.504	240	789134	20.000	ng/ul	0.00
88) Perylene-d12	24.762	264	944712	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	6449	2.619	ng/uL	0.00
4) Pyridine-d5	3.593	84	98047	14.502	ng/ul	0.00
7) Phenol-d5	6.828	99	140219	17.406	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	81621	17.238	ng/ul	0.00
11) 2-Chlorophenol-d4	7.181	132	105201	18.404	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	120277	18.708	ng/ul	0.00
21) Nitrobenzene-d5	8.775	128	53691	18.898	ng/ul	0.00
24) 2-Nitrophenol-d4	9.487	143	65215	19.826	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.028	165	132702	18.982	ng/ul	0.00
31) 4-Chloroaniline-d4	10.534	131	150343	18.200	ng/ul	0.00
46) Dimethylphthalate-d6	13.657	166	484342	20.610	ng/ul	0.00
49) Acenaphthylene-d8	13.939	160	514632	20.619	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.498	143	59799m	15.168	ng/ul	-0.01
60) Fluorene-d10	15.263	176	428469	21.021	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.422	200	47651	9.923	ng/ul	0.02
73) Anthracene-d10	17.169	188	738665	21.504	ng/ul	0.00
81) Pyrene-d10	19.545	212	921274	22.077	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.539	264	1095128	21.706	ng/ul	0.02

Target Compounds Qvalue

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

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