

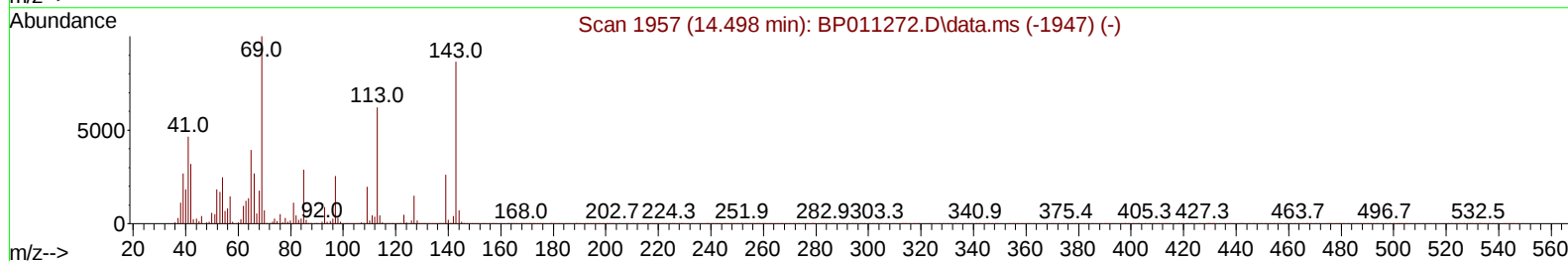
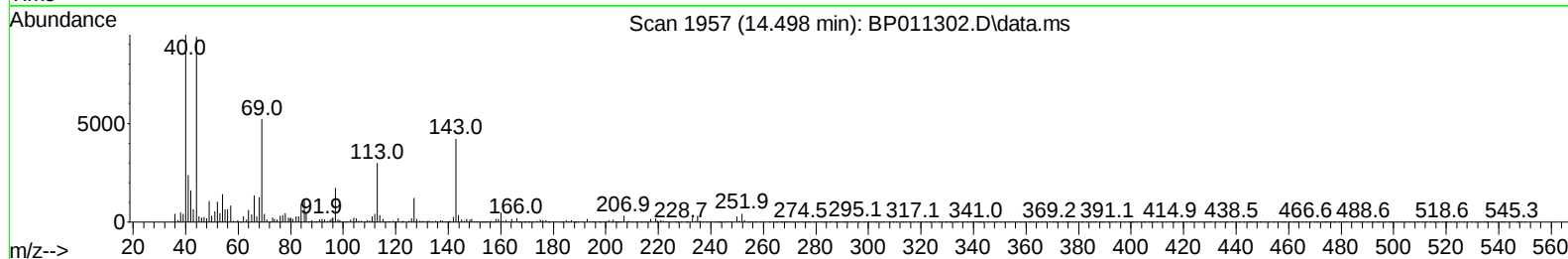
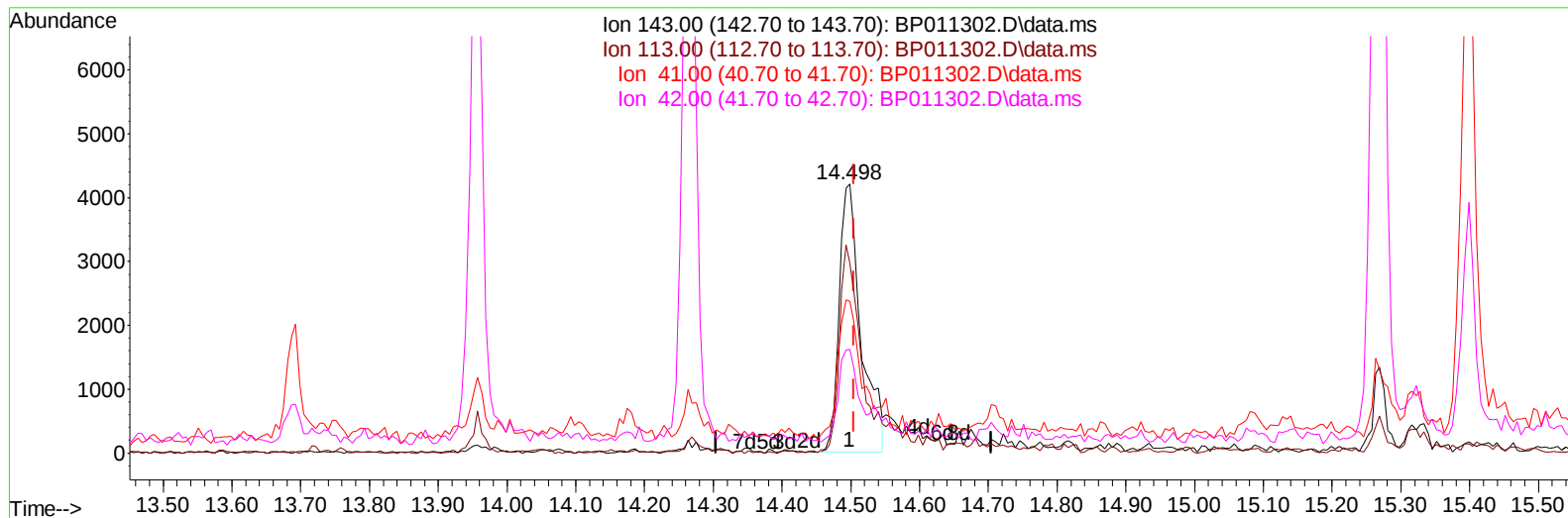
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
Data File : BP011302.D
Acq On : 06 Aug 2022 06:06
Operator : CG/JU
Sample : N4026-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBJJ3

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 08/08/2022
Supervised By :Sohil Jodhani 08/10/2022

Quant Time: Aug 06 06:48:30 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP080222.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Aug 02 23:52:37 2022
Response via : Initial Calibration



TIC: BP011302.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.006) 2.36 ng/u1

response 9319

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	70.72
41.00	54.40	56.57
42.00	35.50	38.57

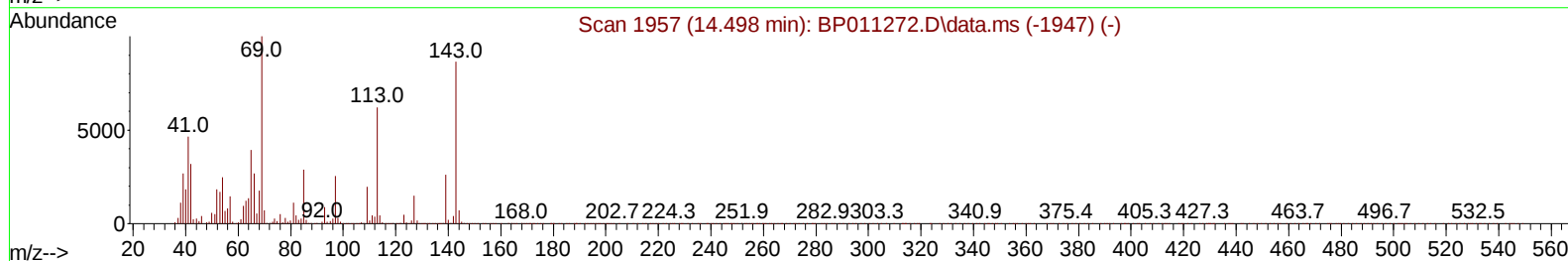
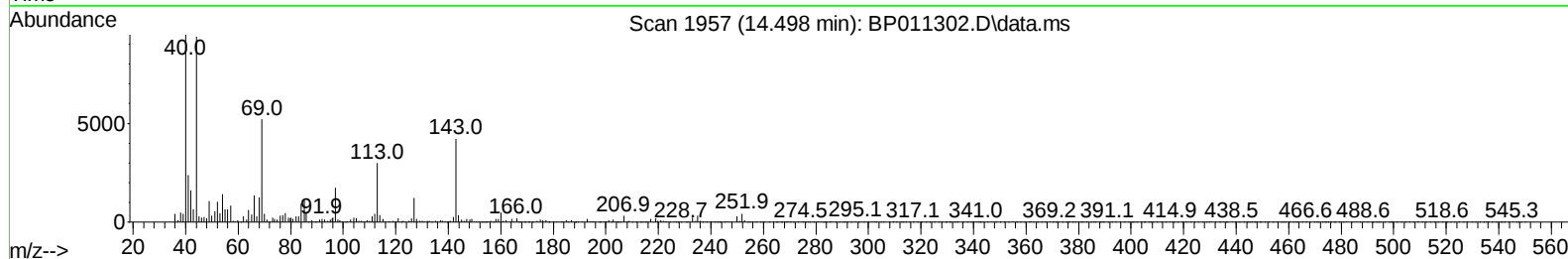
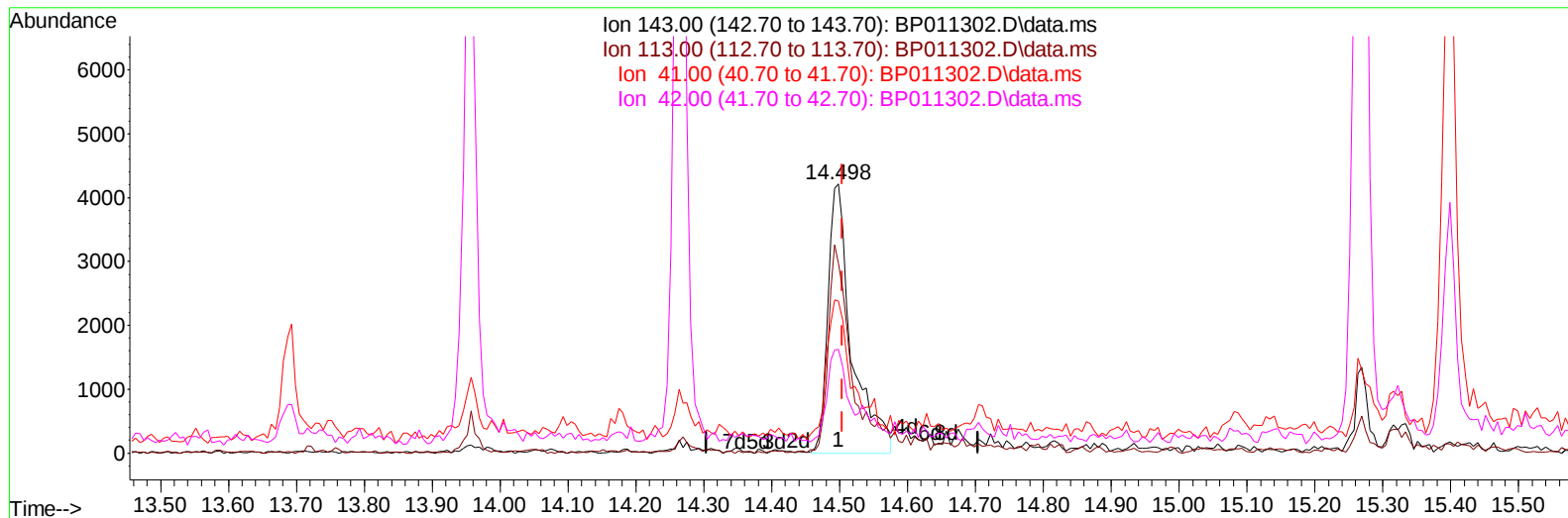
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TIC: BP011302.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.006) 2.59 ng/ul m

response 10256

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	70.72
41.00	54.40	56.57
42.00	35.50	38.57

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	105415	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.381	136	477524	20.000	ng/ul	-0.02
38) Acenaphthene-d10	14.263	164	339727	20.000	ng/ul	#-0.01
64) Phenanthrene-d10	17.034	188	772991	20.000	ng/ul	-0.01
79) Chrysene-d12	21.157	240	933089	20.000	ng/ul	-0.02
88) Perylene-d12	23.469	264	883521	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	2582	1.980	ng/uL	0.00
4) Pyridine-d5	3.505	84	21563	6.132	ng/ul	0.00
7) Phenol-d5	6.775	99	22208	2.528	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	6.940	67	149914	23.649	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.128	132	92771	12.956	ng/ul	-0.01
15) 4-Methylphenol-d8	8.316	113	43211	5.978	ng/ul	-0.01
21) Nitrobenzene-d5	8.758	128	117951	34.341	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.475	143	75943	23.588	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.010	165	109141	17.777	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.534	131	57118	6.014	ng/ul	-0.01
46) Dimethylphthalate-d6	13.693	166	492557	21.169	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	737312	25.186	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.498	143	10256m	2.592	ng/ul	0.00
60) Fluorene-d10	15.269	176	696578	35.183	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.398	200	90229	25.138	ng/ul	-0.02
73) Anthracene-d10	17.134	188	1380723	40.643	ng/ul	-0.01
81) Pyrene-d10	19.392	212	1761069	37.329	ng/ul	-0.02
92) Benzo(a)pyrene-d12	23.322	264	1556525	35.354	ng/ul	-0.02

Target Compounds Qvalue

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

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