

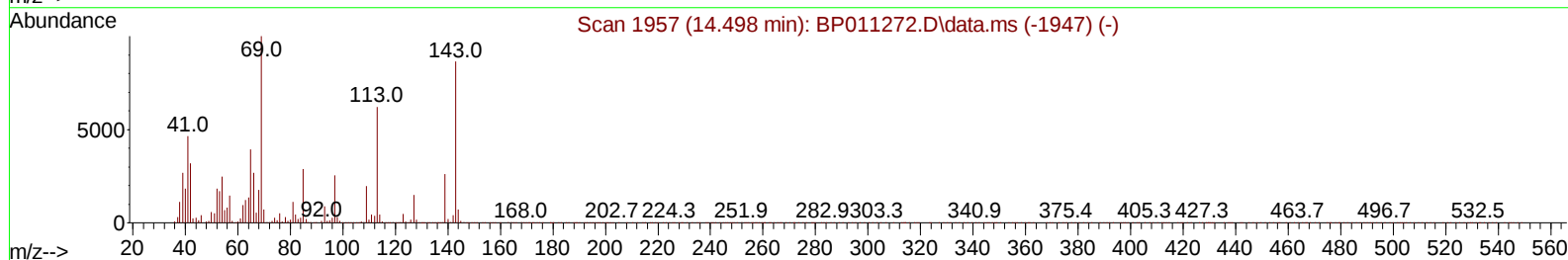
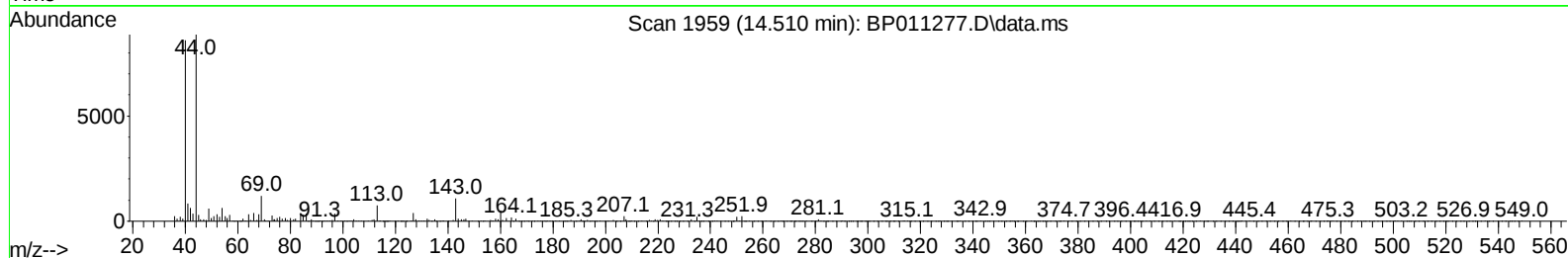
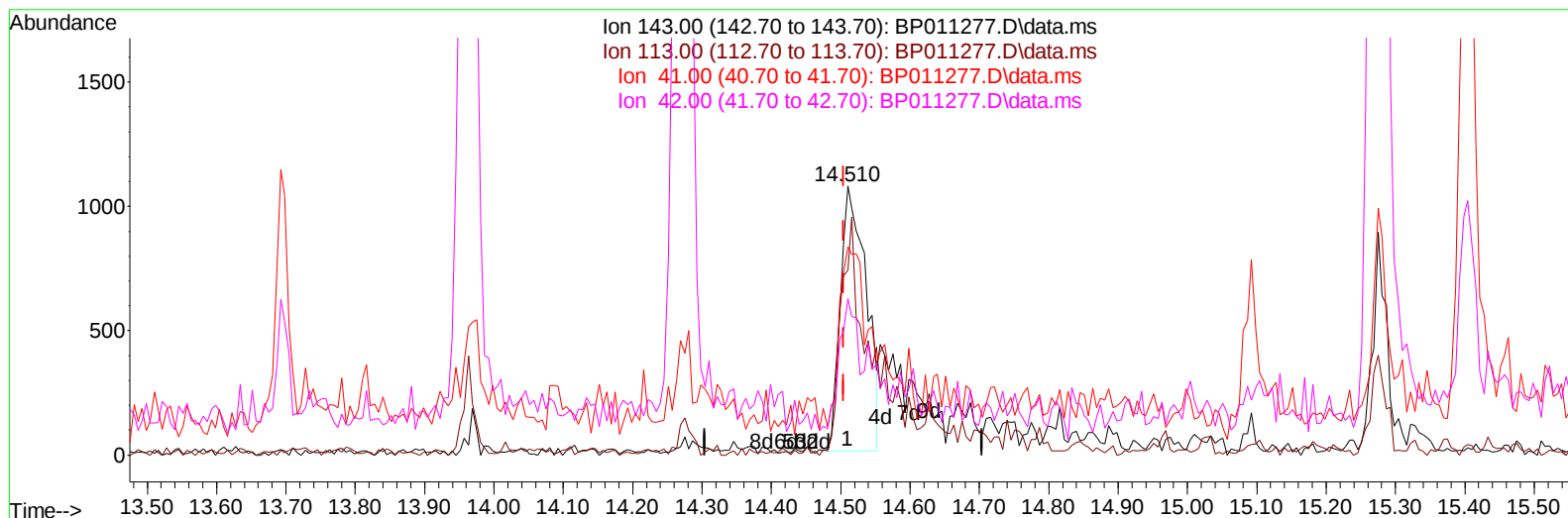
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
Data File : BP011277.D
Acq On : 05 Aug 2022 13:42
Operator : CG/JU
Sample : N4026-03
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBJG8

Manual IntegrationsAPPROVED

Quant Time: Aug 05 23:01:49 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

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



TIC: BP011277.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.006) 1.14 ng/u1

response 2710

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	68.83
41.00	54.40	77.52#
42.00	35.50	58.09#

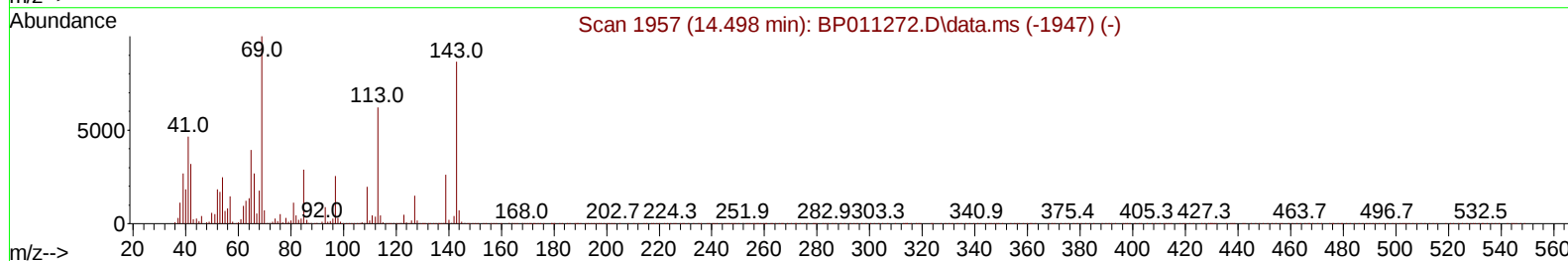
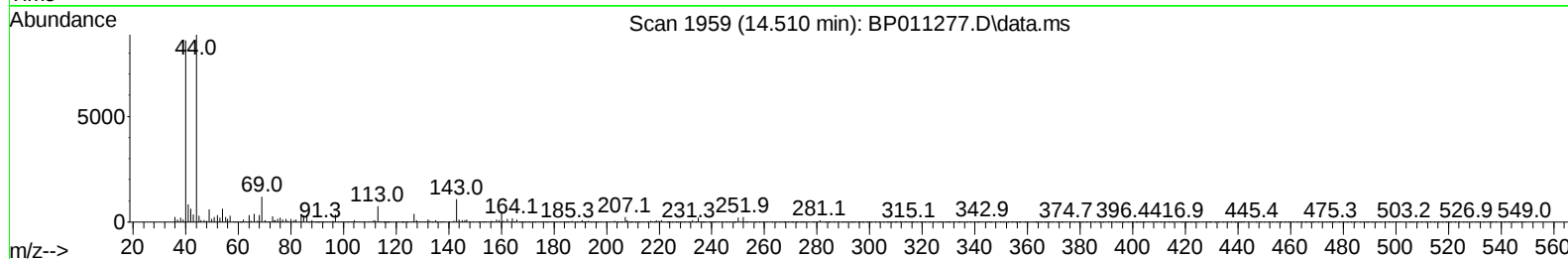
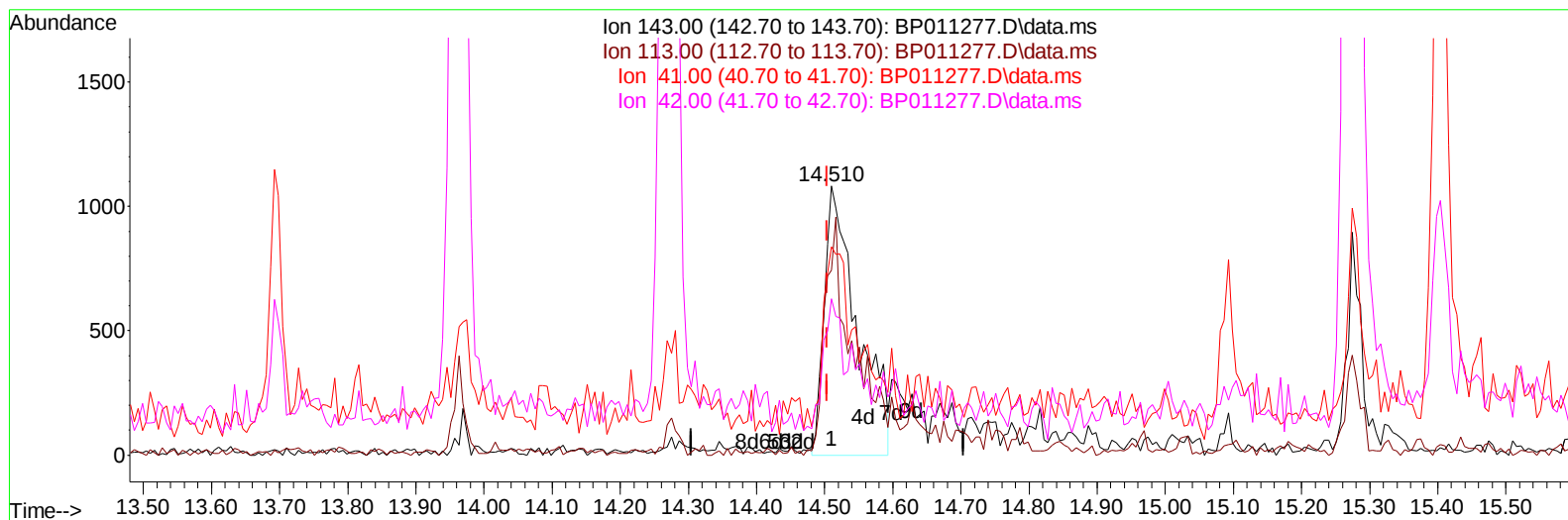
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(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.006) 1.54 ng/ul m

response 3646

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	68.83
41.00	54.40	77.52#
42.00	35.50	58.09#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	91127	20.000	ng/ul	0.00
20) Naphthalene-d8	10.393	136	343393	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	203632	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.045	188	421000	20.000	ng/ul	0.00
79) Chrysene-d12	21.169	240	595290	20.000	ng/ul	0.00
88) Perylene-d12	23.486	264	642570	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	2917	2.588	ng/uL	0.00
4) Pyridine-d5	3.511	84	18835	6.196	ng/ul	0.00
7) Phenol-d5	6.781	99	14712	1.937	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	152365	27.805	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	48826	7.888	ng/ul	0.00
15) 4-Methylphenol-d8	8.322	113	24937	3.991	ng/ul	0.00
21) Nitrobenzene-d5	8.763	128	103481	41.896	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	32275	13.940	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.022	165	42232	9.566	ng/ul	0.00
31) 4-Chloroaniline-d4	10.546	131	46134	6.754	ng/ul	0.00
46) Dimethylphthalate-d6	13.698	166	262281	18.806	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	526485	30.003	ng/ul	0.00
54) 4-Nitrophenol-d4	14.510	143	3646m	1.537	ng/ul	0.00
60) Fluorene-d10	15.275	176	493318	41.570	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	23657	12.102	ng/ul	-0.01
73) Anthracene-d10	17.139	188	1029838	55.659	ng/ul	0.00
81) Pyrene-d10	19.404	212	1146153	38.081	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.339	264	1195927	37.349	ng/ul	0.00

Target Compounds Qvalue

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

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