

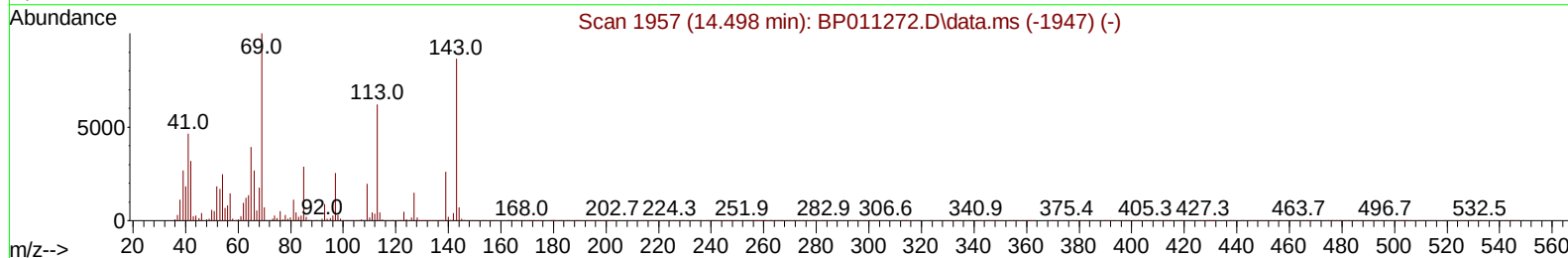
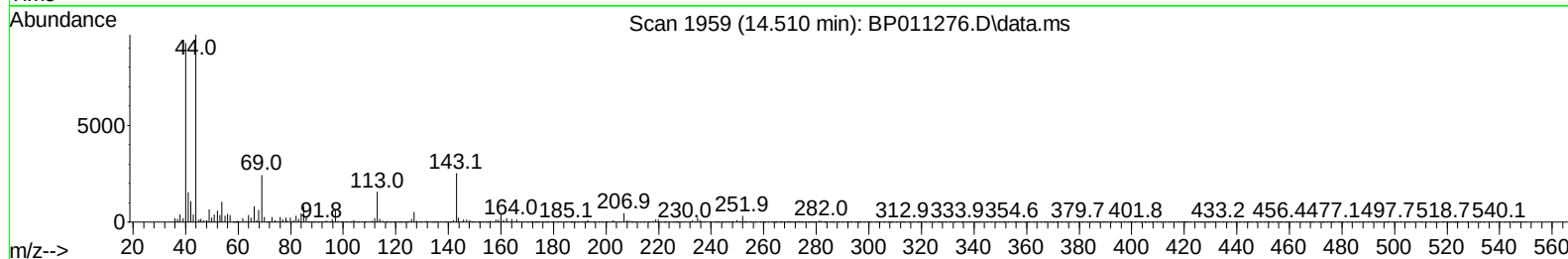
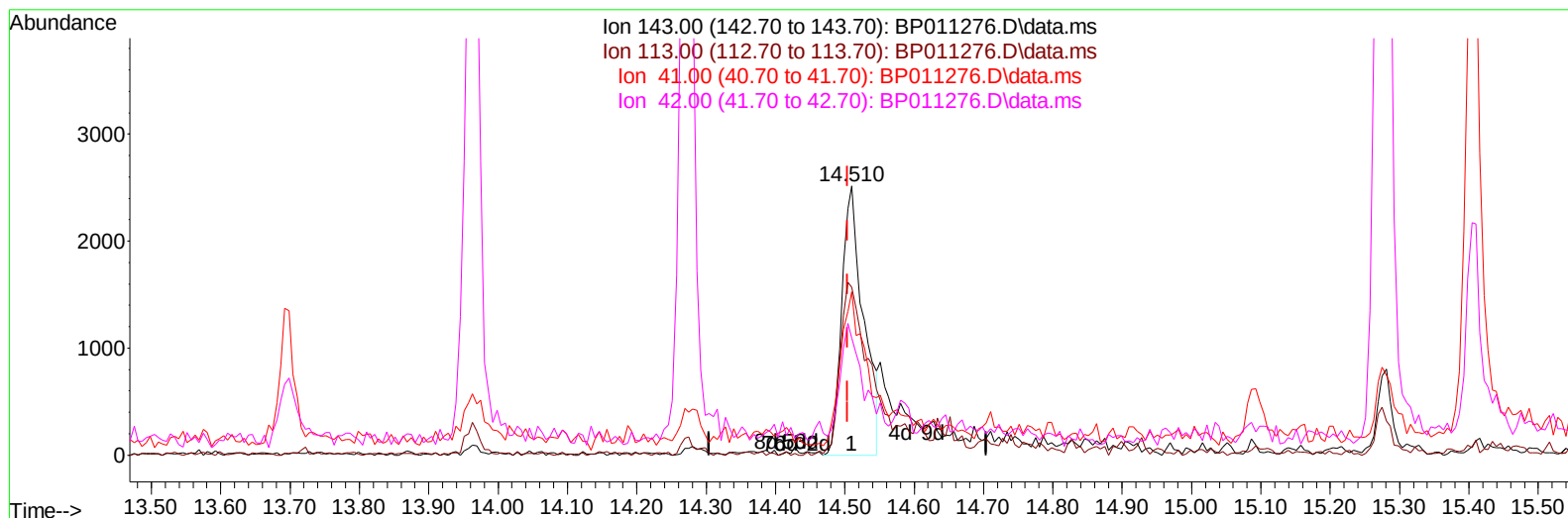
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
Data File : BP011276.D
Acq On : 05 Aug 2022 13:07
Operator : CG/JU
Sample : N4026-01
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GBJG6

Manual IntegrationsAPPROVED

Quant Time: Aug 05 23:01:38 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: BP011276.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.006) 2.23 ng/ul

response 5448

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	62.49
41.00	54.40	60.74
42.00	35.50	43.52#

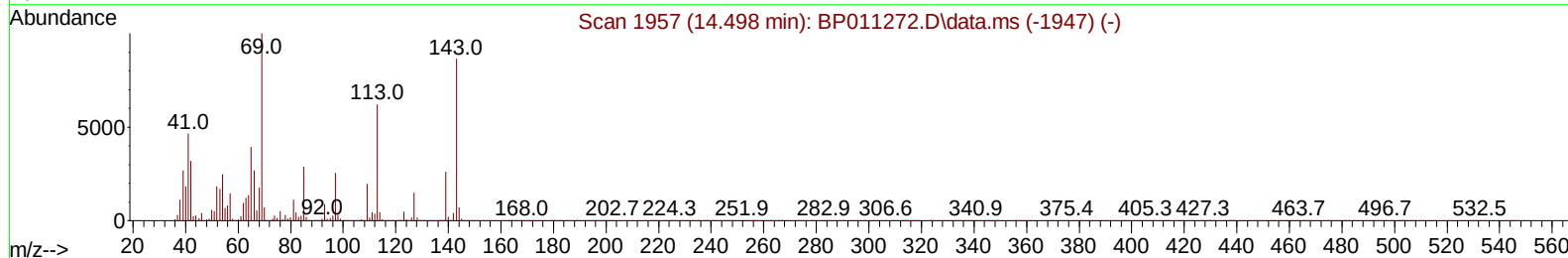
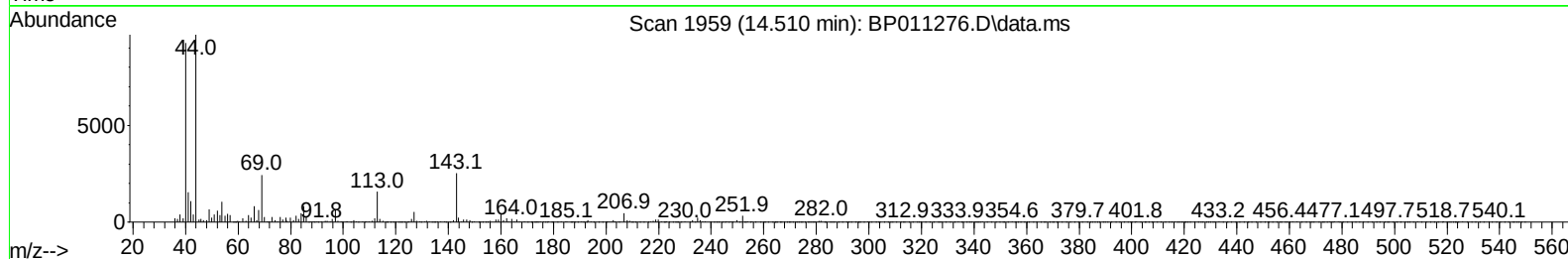
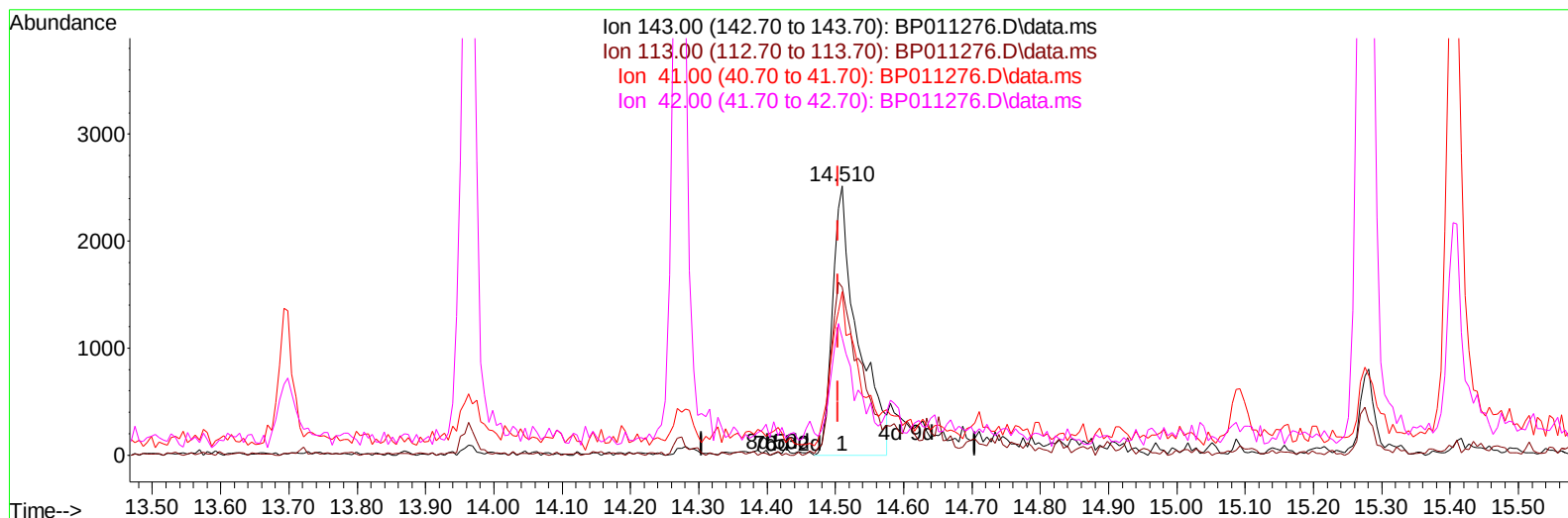
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TIC: BP011276.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.510min (+ 0.006) 2.64 ng/ul m

response 6452

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	62.49
41.00	54.40	60.74
42.00	35.50	43.52#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	91686	20.000	ng/ul	0.00
20) Naphthalene-d8	10.393	136	366203	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.275	164	209690	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.045	188	422805	20.000	ng/ul	0.00
79) Chrysene-d12	21.169	240	589057	20.000	ng/ul	0.00
88) Perylene-d12	23.486	264	637831	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.093	96	3622	3.194	ng/uL	0.00
4) Pyridine-d5	3.511	84	23345	7.632	ng/ul	0.00
7) Phenol-d5	6.781	99	27660	3.620	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	170931	31.002	ng/ul	0.00
11) 2-Chlorophenol-d4	7.134	132	103930	16.688	ng/ul	0.00
15) 4-Methylphenol-d8	8.322	113	53982	8.586	ng/ul	0.00
21) Nitrobenzene-d5	8.763	128	111115	42.185	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	68713	27.830	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.016	165	100384	21.321	ng/ul	0.00
31) 4-Chloroaniline-d4	10.546	131	50467	6.929	ng/ul	0.00
46) Dimethylphthalate-d6	13.698	166	379578	26.430	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	579034	32.045	ng/ul	0.00
54) 4-Nitrophenol-d4	14.510	143	6452m	2.642	ng/ul	0.00
60) Fluorene-d10	15.275	176	530732	43.431	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	52785	26.887	ng/ul	-0.01
73) Anthracene-d10	17.145	188	1113915	59.946	ng/ul	0.00
81) Pyrene-d10	19.404	212	1236329	41.511	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.339	264	1308437	41.166	ng/ul	0.00

Target Compounds Qvalue

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

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