

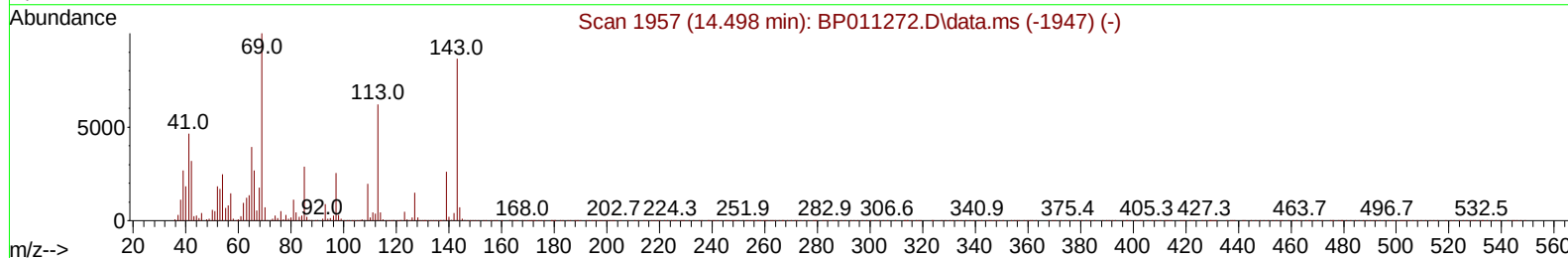
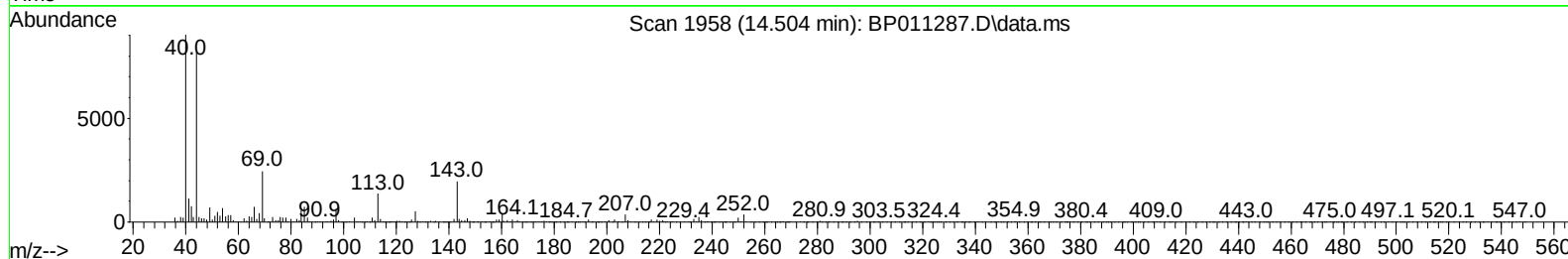
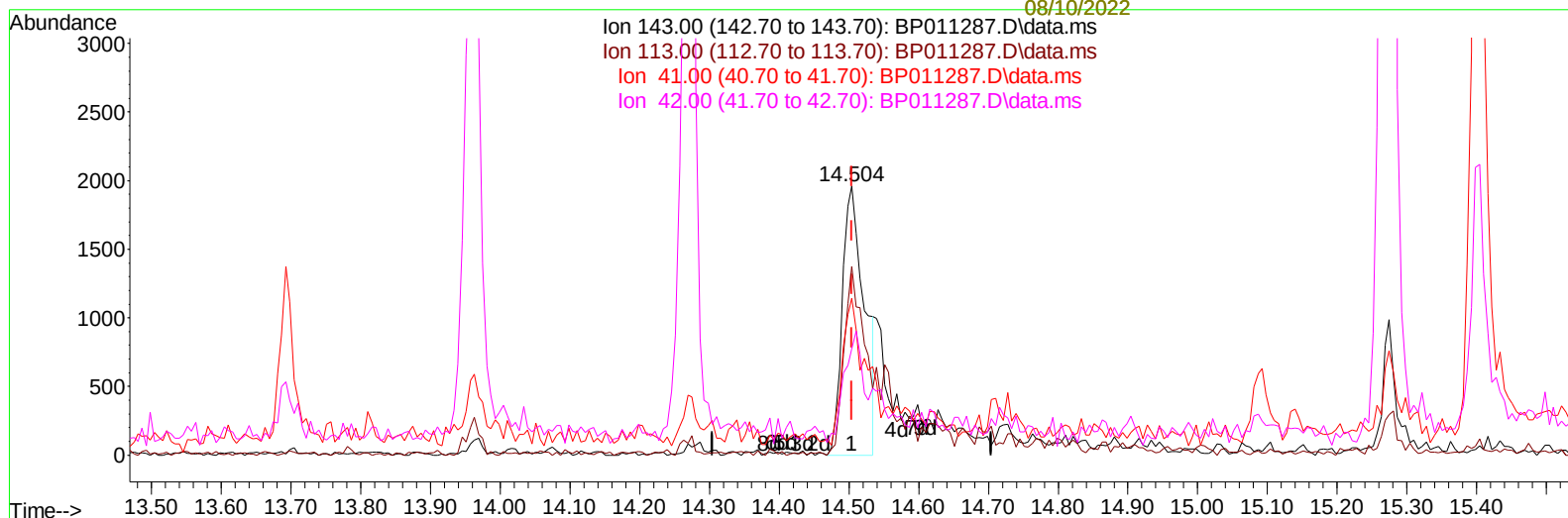
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
 Data File : BP011287.D
 Acq On : 05 Aug 2022 20:07
 Operator : CG/JU
 Sample : N4026-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJH2

Manual IntegrationsAPPROVED

Quant Time: Aug 05 23:03:50 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: BP011287.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (-0.000) 2.09 ng/u1

response 4262

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	70.15
41.00	54.40	58.42
42.00	35.50	39.49

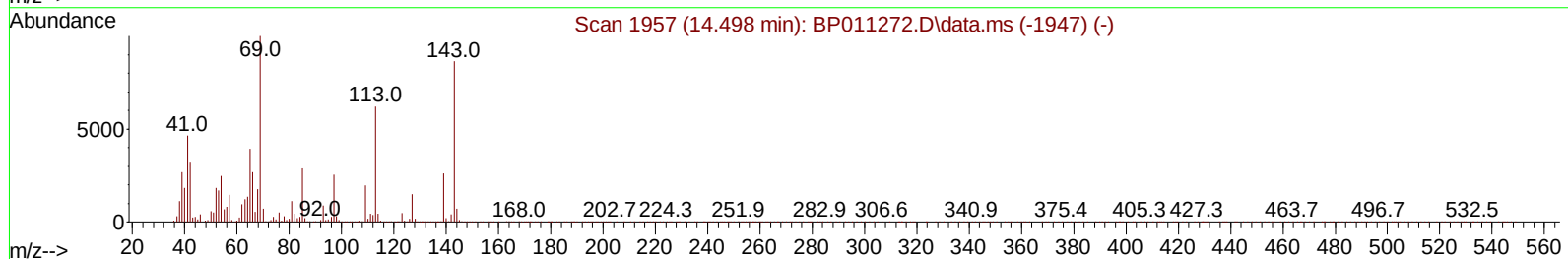
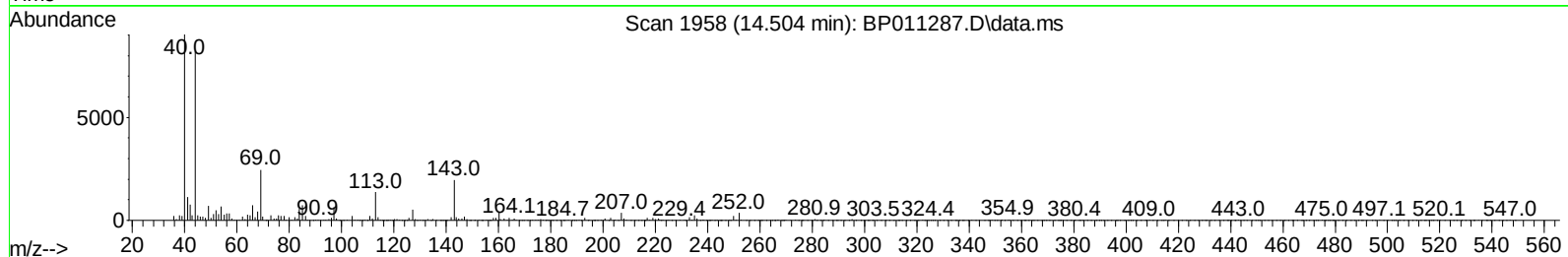
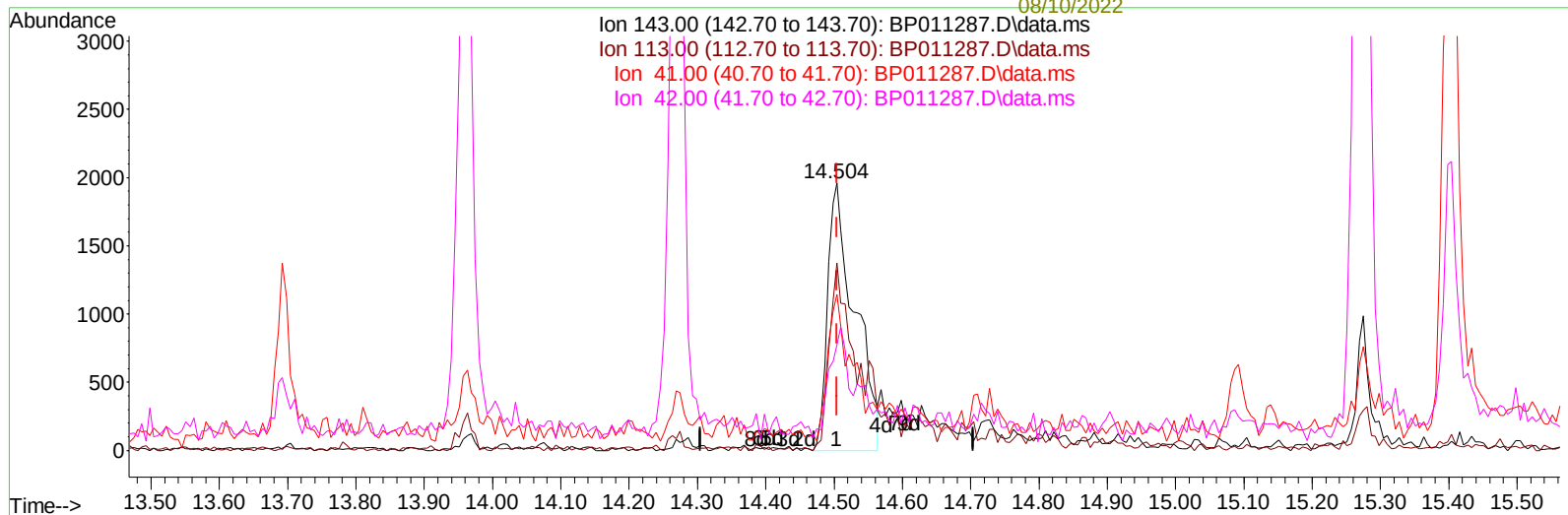
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TIC: BP011287.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (-0.000) 2.63 ng/ul m

response 5373

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	70.15
41.00	54.40	58.42
42.00	35.50	39.49

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.599	152		90792	20.000	ng/ul	0.00
20) Naphthalene-d8	10.387	136		337247	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.269	164		175100	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.039	188		327008	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240		488537	20.000	ng/ul	-0.01
88) Perylene-d12	23.474	264		542017	20.000	ng/ul	-0.02
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.087	96		3174	2.827	ng/uL	0.00
4) Pyridine-d5	3.511	84		12503	4.128	ng/ul	0.00
7) Phenol-d5	6.781	99		22959	3.034	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67		157907	28.922	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132		92891	15.062	ng/ul	-0.01
15) 4-Methylphenol-d8	8.316	113		45509	7.310	ng/ul	-0.01
21) Nitrobenzene-d5	8.763	128		114368	47.148	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143		67563	29.714	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.016	165		92107	21.243	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131		41194	6.141	ng/ul	0.00
46) Dimethylphthalate-d6	13.692	166		321517	26.810	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160		513372	34.023	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143		5373m	2.635	ng/ul	0.00
60) Fluorene-d10	15.275	176		449338	44.034	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200		52006	34.250	ng/ul	-0.01
73) Anthracene-d10	17.139	188		900658	62.669	ng/ul	0.00
81) Pyrene-d10	19.398	212		1028963	41.657	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.321	264		1123878	41.610	ng/ul	-0.02

Target Compounds Qvalue

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

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