

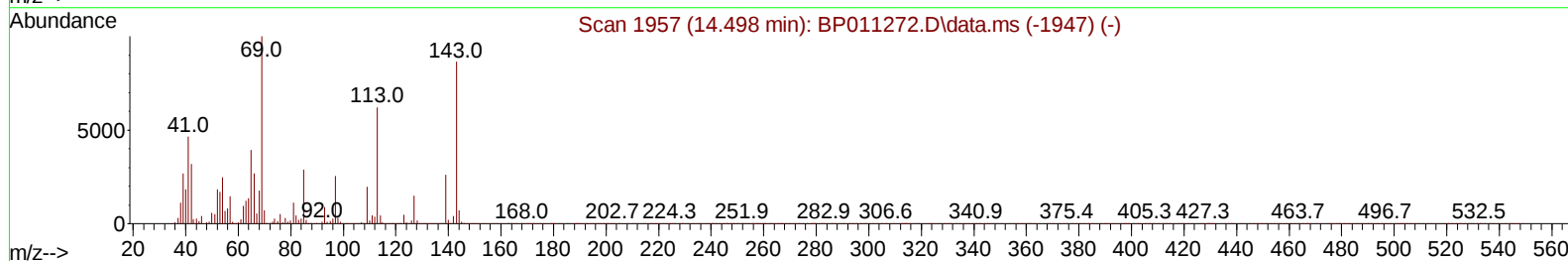
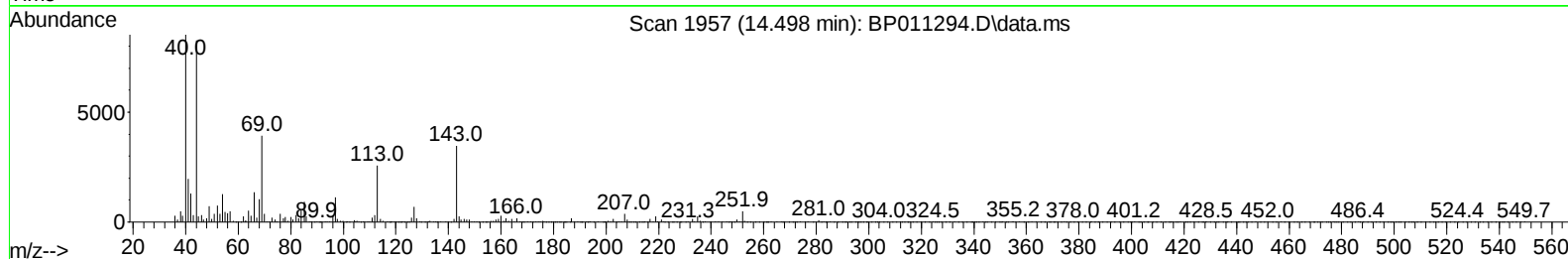
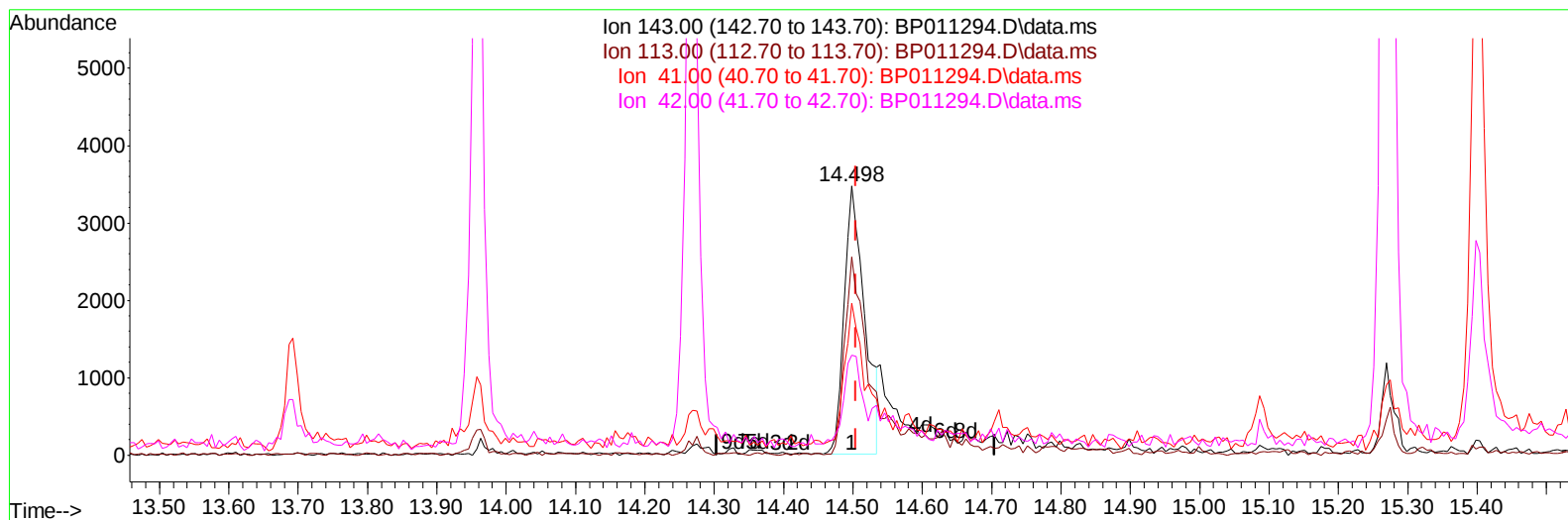
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
 Data File : BP011294.D
 Acq On : 06 Aug 2022 00:10
 Operator : CG/JU
 Sample : N4026-12
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJH6

Manual IntegrationsAPPROVED

Quant Time: Aug 06 00:59:09 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: BP011294.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.006) 2.44 ng/u1

response 7051

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	73.74
41.00	54.40	56.46
42.00	35.50	37.19

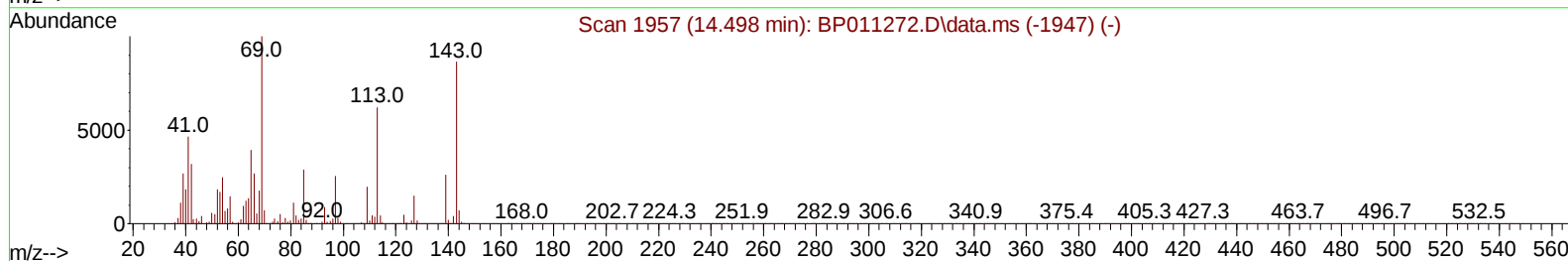
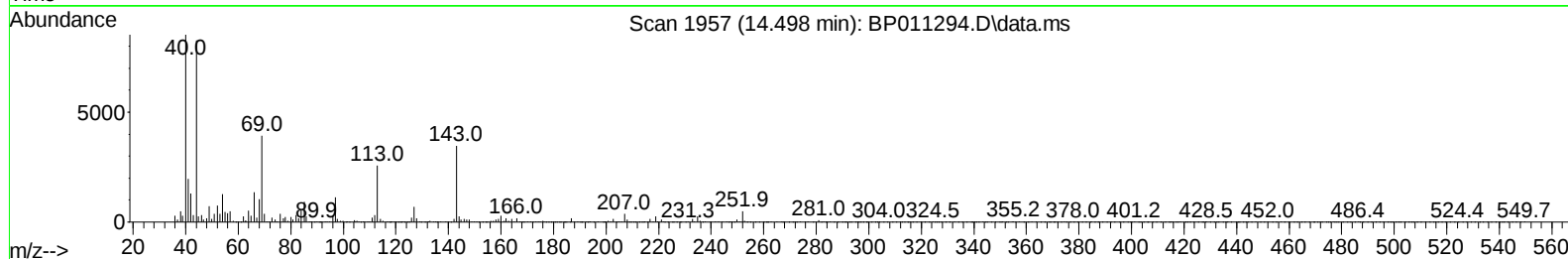
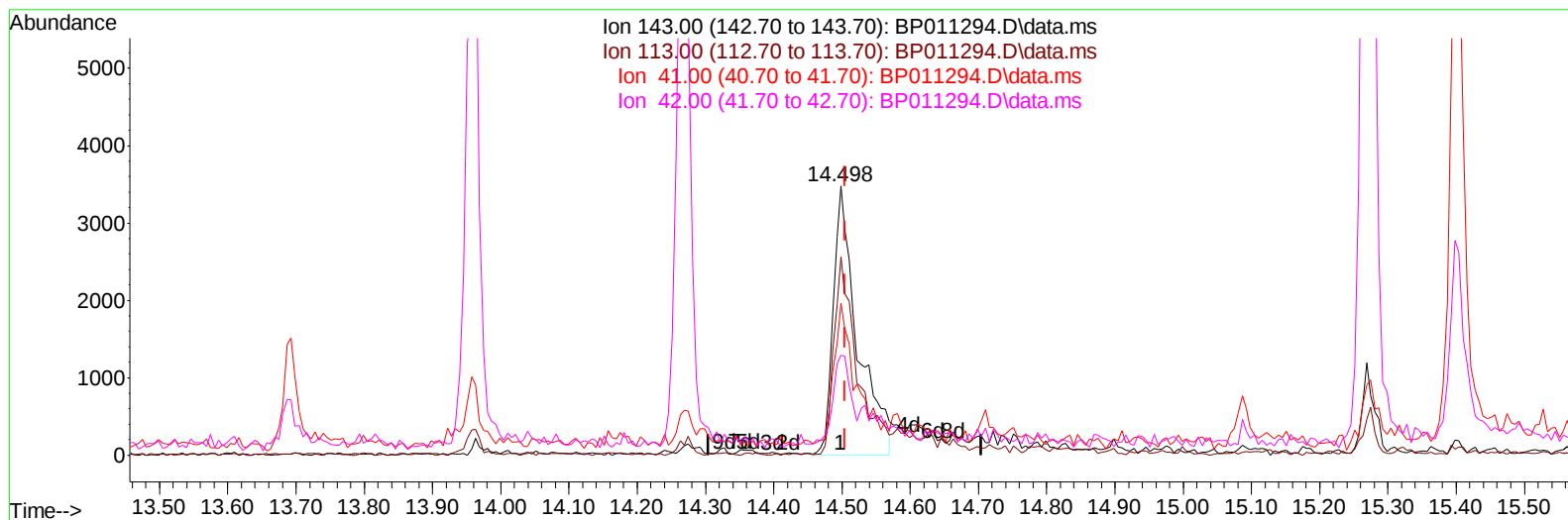
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TIC: BP011294.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.498min (-0.006) 2.98 ng/ul m

response 8610

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	73.74
41.00	54.40	56.46
42.00	35.50	37.19

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	103938	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.387	136	418314	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.269	164	247836	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.039	188	511916	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	702870	20.000	ng/ul	-0.01
88) Perylene-d12	23.474	264	715454	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	2778	2.161	ng/uL	0.00
4) Pyridine-d5	3.505	84	21359	6.160	ng/ul	0.00
7) Phenol-d5	6.781	99	25532	2.948	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	162965	26.073	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.128	132	103983	14.728	ng/ul	-0.01
15) 4-Methylphenol-d8	8.316	113	53185	7.462	ng/ul	-0.01
21) Nitrobenzene-d5	8.758	128	123190	40.943	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.475	143	76361	27.075	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.016	165	103401	19.226	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	55139	6.627	ng/ul	0.00
46) Dimethylphthalate-d6	13.693	166	427381	25.178	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	661502	30.974	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.498	143	8610m	2.983	ng/ul	0.00
60) Fluorene-d10	15.269	176	612774	42.426	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.404	200	77618	32.653	ng/ul	-0.01
73) Anthracene-d10	17.134	188	1281285	56.950	ng/ul	-0.01
81) Pyrene-d10	19.398	212	1539485	43.320	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.321	264	1473246	41.323	ng/ul	-0.02

Target Compounds Qvalue

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

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