

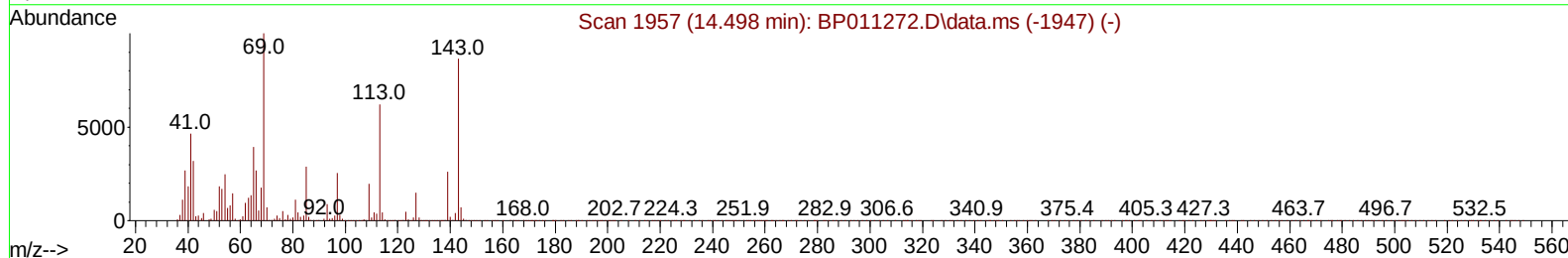
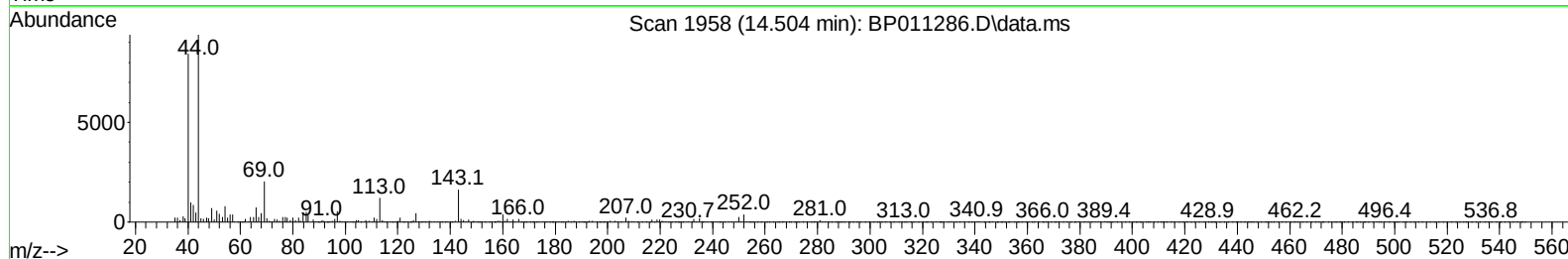
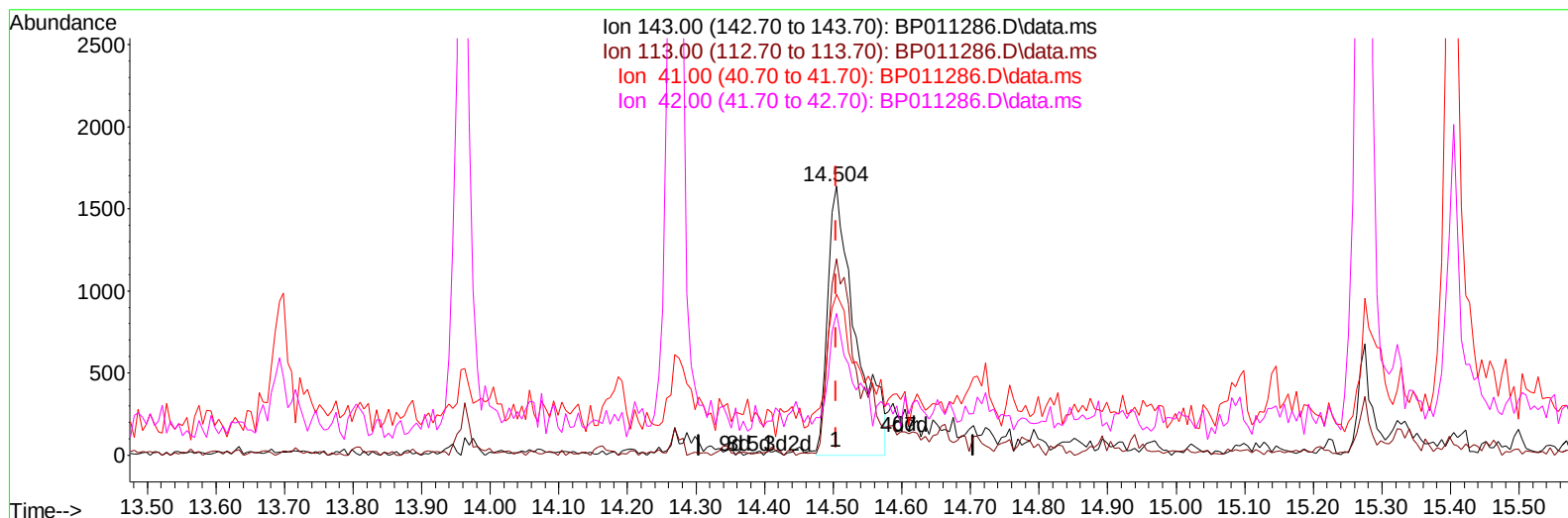
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
 Data File : BP011286.D
 Acq On : 05 Aug 2022 19:32
 Operator : CG/JU
 Sample : N4026-17
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJJ1

Manual IntegrationsAPPROVED

Quant Time: Aug 05 23:03:37 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: BP011286.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (+ 0.000) 2.22 ng/ul m

response 4597

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	72.91
41.00	54.40	59.67
42.00	35.50	52.65#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	94096	20.000	ng/ul	0.00
20) Naphthalene-d8	10.387	136	334954	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.269	164	177964	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.040	188	387070	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	572971	20.000	ng/ul	-0.01
88) Perylene-d12	23.474	264	591183	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	2802	2.408	ng/uL	0.00
4) Pyridine-d5	3.505	84	24260	7.728	ng/ul	0.00
7) Phenol-d5	6.781	99	15320	1.954	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.946	67	115323	20.381	ng/ul	0.00
11) 2-Chlorophenol-d4	7.128	132	68151	10.662	ng/ul	-0.01
15) 4-Methylphenol-d8	8.316	113	26743	4.145	ng/ul	-0.01
21) Nitrobenzene-d5	8.764	128	78664	32.651	ng/ul	0.00
24) 2-Nitrophenol-d4	9.481	143	46814	20.730	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.016	165	62723	14.565	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	28310	4.249	ng/ul	0.00
46) Dimethylphthalate-d6	13.693	166	229867	18.859	ng/ul	0.00
49) Acenaphthylene-d8	13.963	160	311473	20.310	ng/ul	0.00
54) 4-Nitrophenol-d4	14.504	143	4597m	2.218	ng/ul	0.00
60) Fluorene-d10	15.275	176	301726	29.092	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	37900	21.087	ng/ul	-0.01
73) Anthracene-d10	17.140	188	610256	35.873	ng/ul	0.00
81) Pyrene-d10	19.398	212	820755	28.332	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.327	264	851308	28.898	ng/ul	-0.02

Target Compounds Qvalue

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

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