

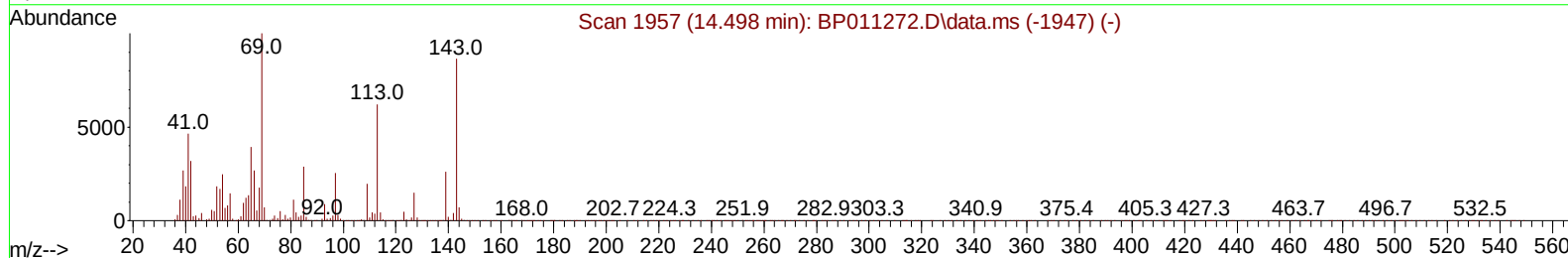
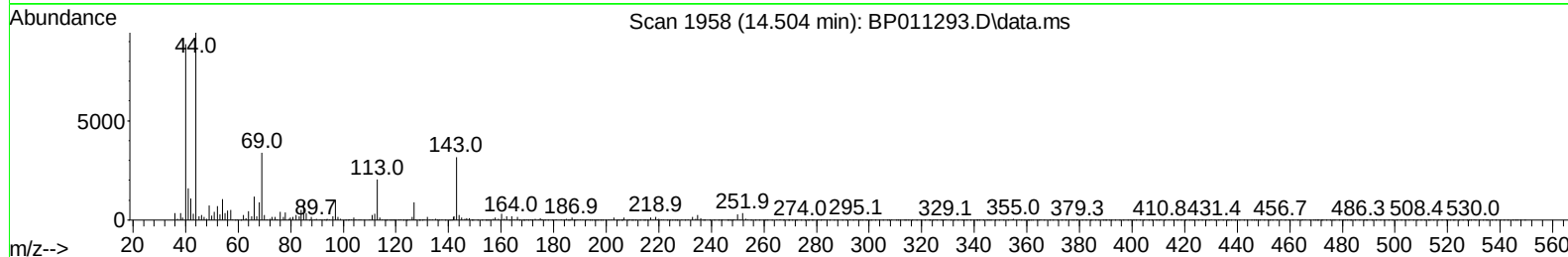
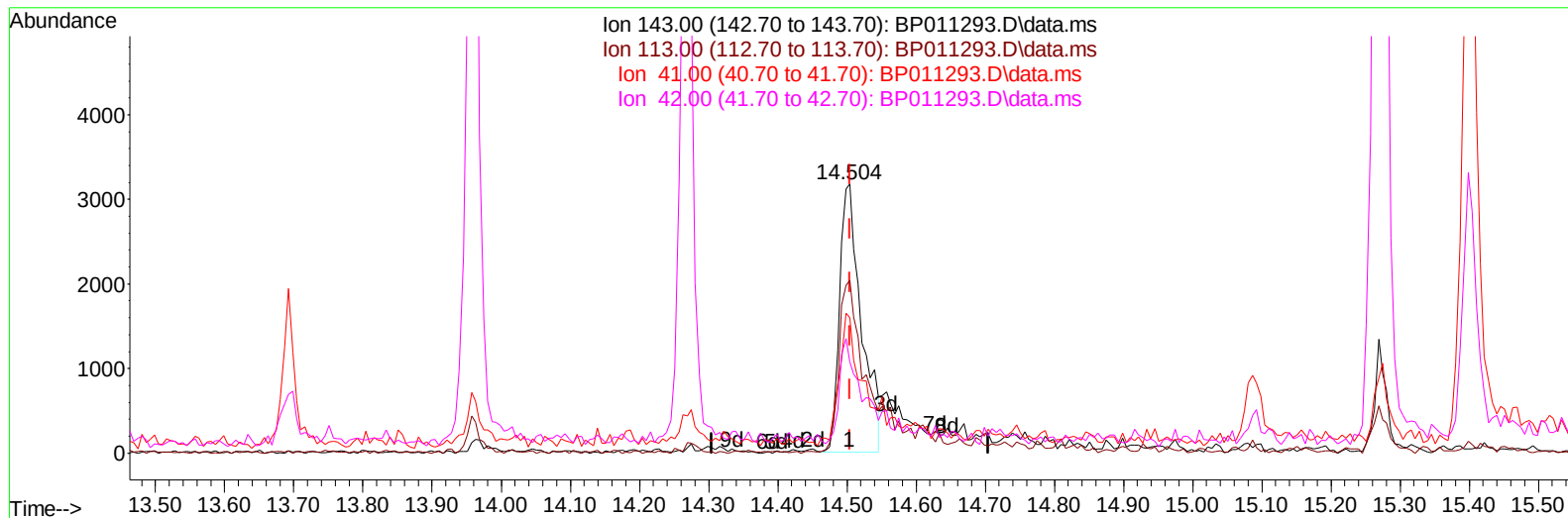
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080522\
 Data File : BP011293.D
 Acq On : 05 Aug 2022 23:36
 Operator : CG/JU
 Sample : N4026-15
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GBJH9

Manual IntegrationsAPPROVED

Quant Time: Aug 06 00:14:42 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: BP011293.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (-0.000) 2.88 ng/u1

response 6965

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	64.31
41.00	54.40	50.25
42.00	35.50	33.81

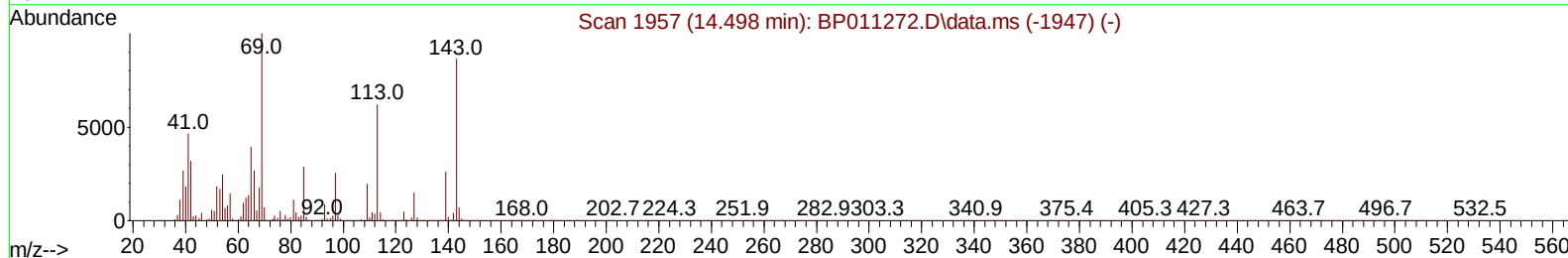
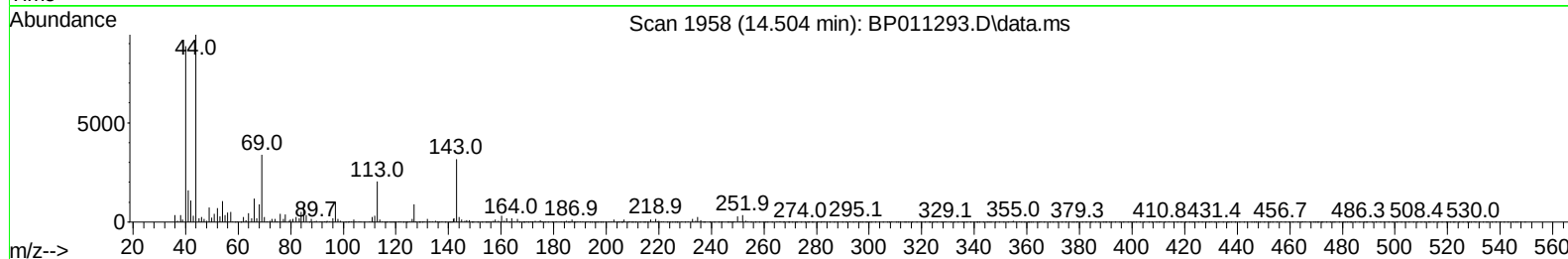
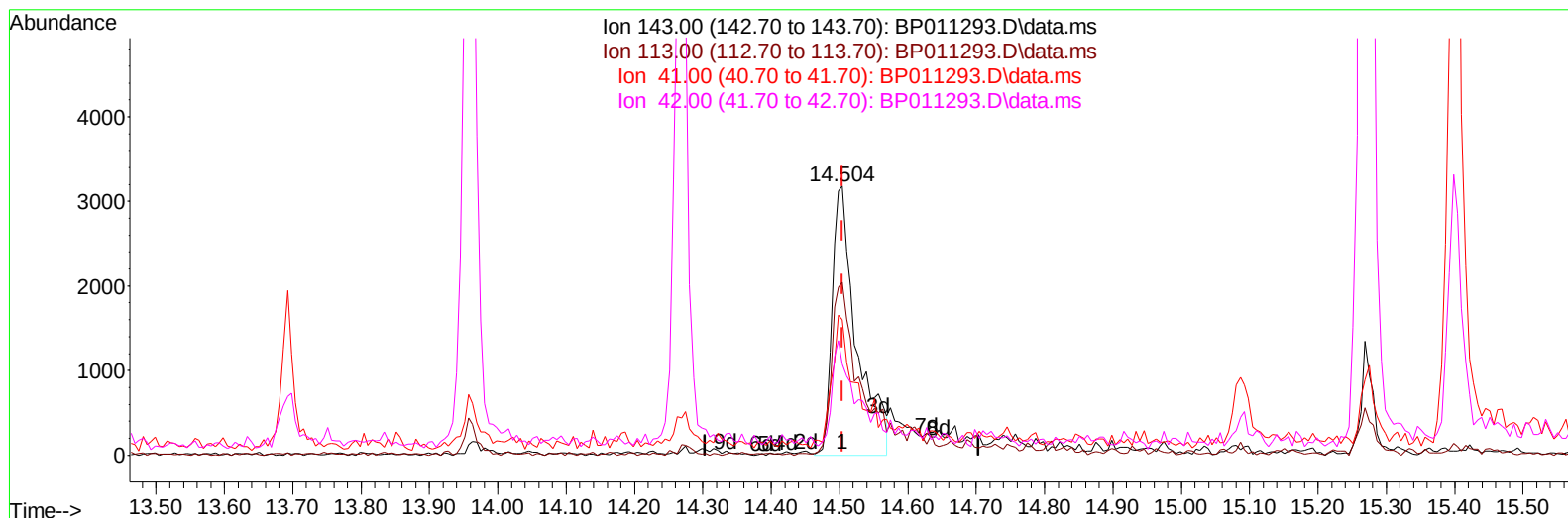
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TIC: BP011293.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.504min (-0.000) 3.27 ng/ul m

response 7900

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.70	64.31
41.00	54.40	50.25
42.00	35.50	33.81

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.593	152	93165	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.387	136	356874	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.269	164	207426	20.000	ng/ul	# 0.00
64) Phenanthrene-d10	17.039	188	417072	20.000	ng/ul	0.00
79) Chrysene-d12	21.163	240	564419	20.000	ng/ul	-0.01
88) Perylene-d12	23.468	264	607818	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.087	96	2923	2.537	ng/uL	0.00
4) Pyridine-d5	3.505	84	24508	7.885	ng/ul	0.00
7) Phenol-d5	6.781	99	26628	3.430	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.940	67	191021	34.096	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.128	132	116389	18.391	ng/ul	-0.01
15) 4-Methylphenol-d8	8.316	113	57994	9.078	ng/ul	-0.01
21) Nitrobenzene-d5	8.757	128	143565	55.929	ng/ul	-0.01
24) 2-Nitrophenol-d4	9.475	143	87691	36.445	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.016	165	122927	26.791	ng/ul	0.00
31) 4-Chloroaniline-d4	10.540	131	115056	16.209	ng/ul	0.00
46) Dimethylphthalate-d6	13.692	166	472288	33.244	ng/ul	0.00
49) Acenaphthylene-d8	13.957	160	764103	42.748	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.504	143	7900m	3.270	ng/ul	0.00
60) Fluorene-d10	15.269	176	699009	57.825	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.404	200	83204	42.963	ng/ul	-0.01
73) Anthracene-d10	17.139	188	1346110	73.438	ng/ul	0.00
81) Pyrene-d10	19.398	212	1522031	53.335	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.321	264	1491205	49.233	ng/ul	-0.02

Target Compounds Qvalue

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

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