

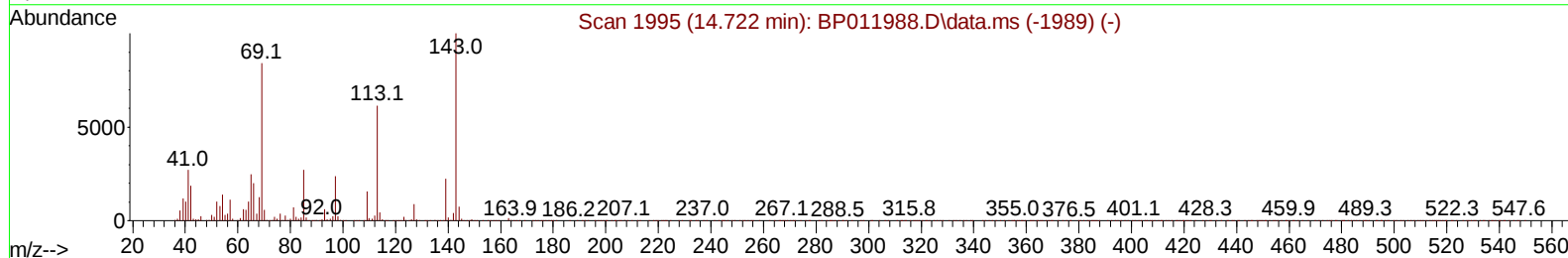
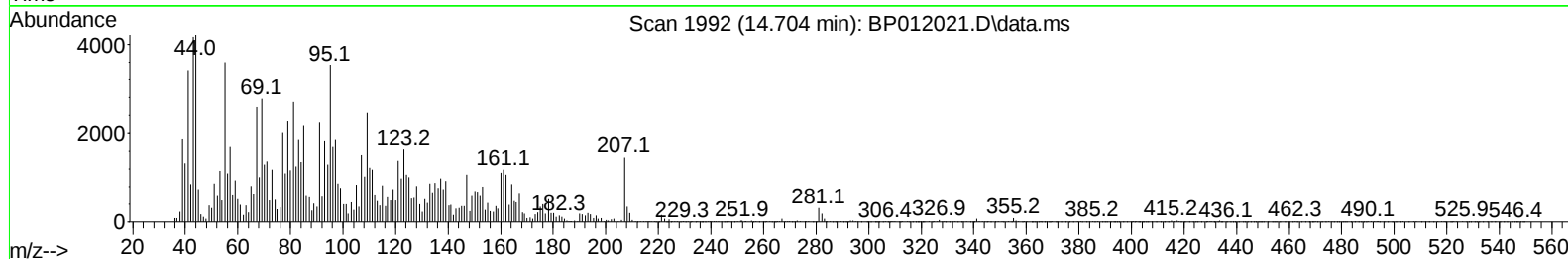
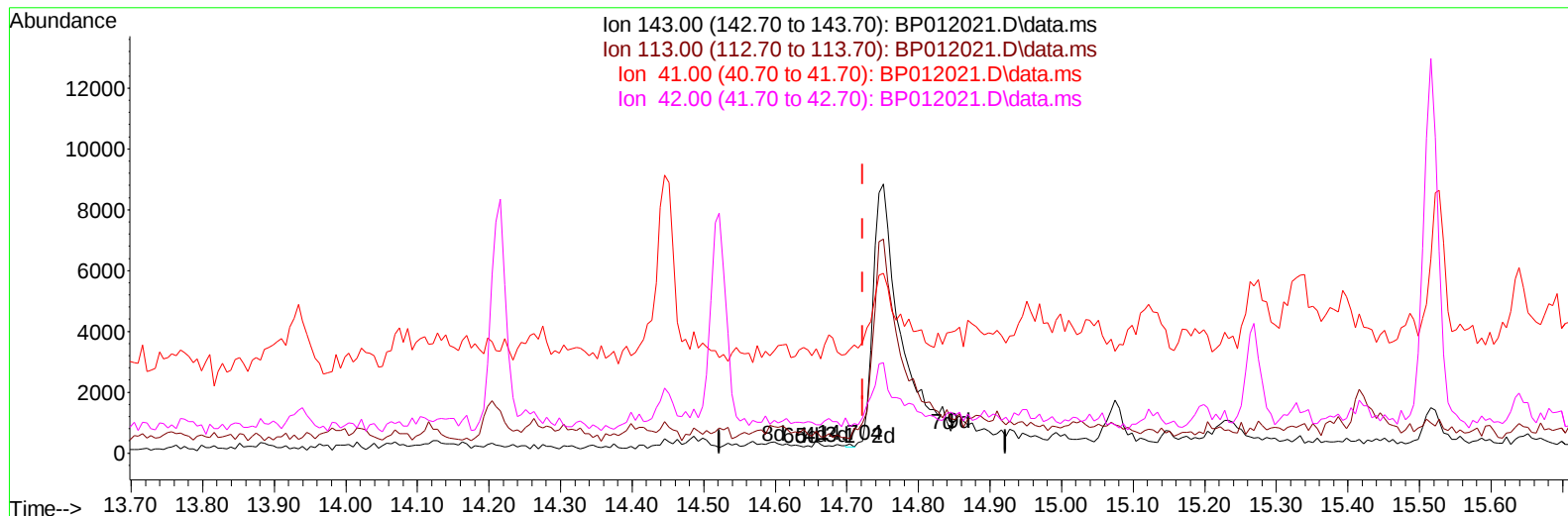
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100522\
Data File : BP012021.D
Acq On : 07 Oct 2022 22:33
Operator : CG/JU
Sample : N4951-03
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
CB8T8

Manual IntegrationsAPPROVED

Quant Time: Oct 07 23:11:44 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP093022.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 07 03:22:44 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 10/10/2022
Supervised By :mohammad ahmed 10/10/2022



TIC: BP012021.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.704min (-0.018) 0.00 ng/u1

response 35

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	162.37#
41.00	24.20	1154.24#
42.00	16.40	293.56#

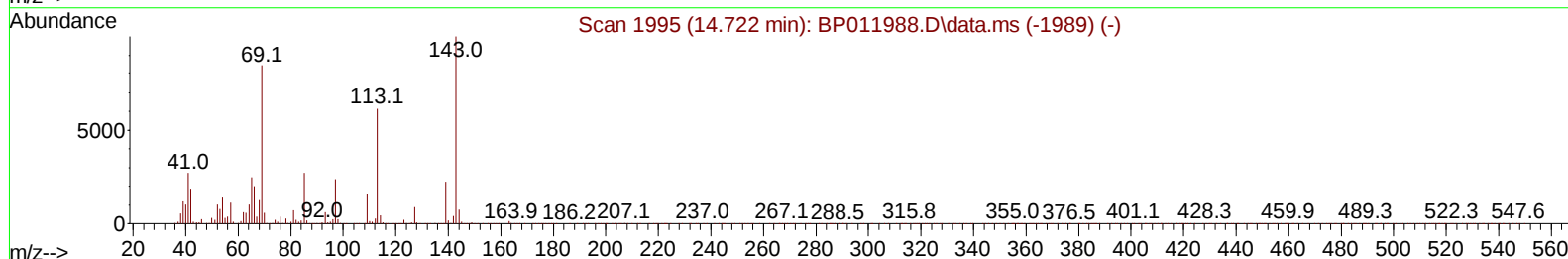
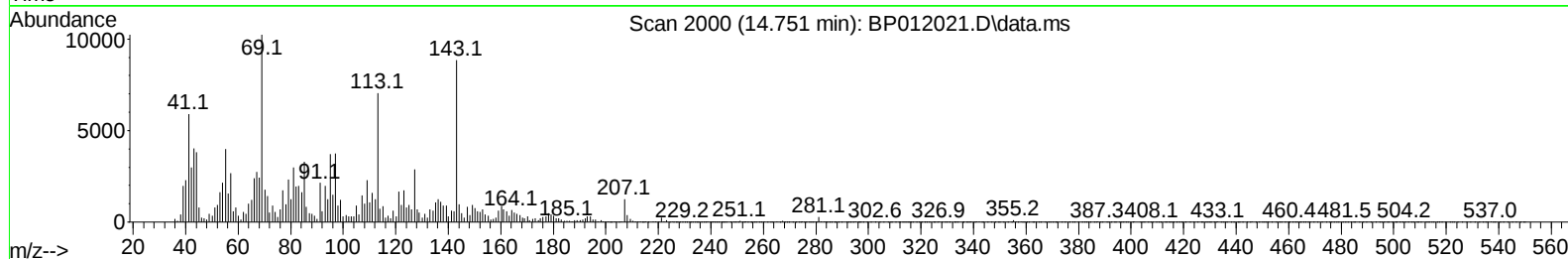
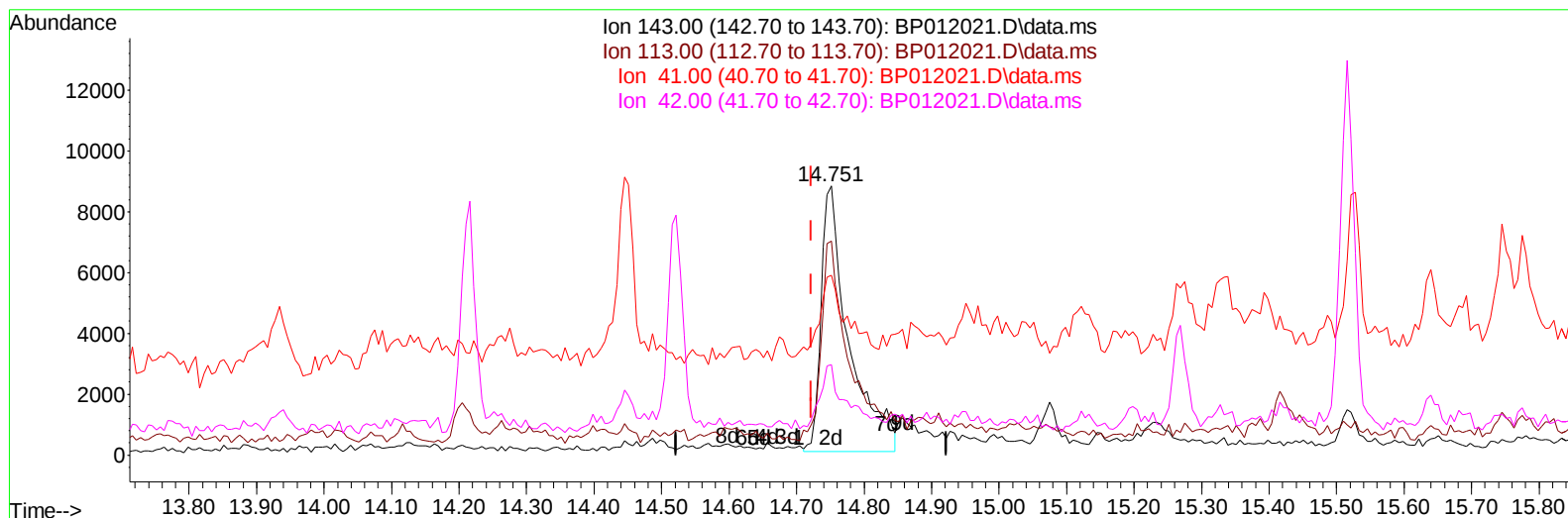
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TIC: BP012021.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (+ 0.029) 3.00 ng/ul m

response 24707

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	59.80	79.64#
41.00	24.20	66.80#
42.00	16.40	33.66#

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 Misc :
 ALS Vial : 17 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.875	152	212435	20.000	ng/ul	0.00
20) Naphthalene-d8	10.681	136	930384	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.522	164	556154	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.280	188	1170809	20.000	ng/ul	0.00
79) Chrysene-d12	21.368	240	1138170	20.000	ng/ul	0.00
88) Perylene-d12	23.798	264	1182243	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.299	96	12785	2.654	ng/uL	0.00
4) Pyridine-d5	3.728	84	159980	11.339	ng/ul	0.01
7) Phenol-d5	7.046	99	116196	6.336	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.205	67	268457	26.409	ng/ul	0.00
11) 2-Chlorophenol-d4	7.405	132	300248	20.739	ng/ul	0.00
15) 4-Methylphenol-d8	8.587	113	212623	14.025	ng/ul	0.00
21) Nitrobenzene-d5	9.046	128	191577	27.034	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	181522	23.941	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.304	165	323707	23.082	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	561079	26.243	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	1193095	29.999	ng/ul	0.00
49) Acenaphthylene-d8	14.216	160	1398363	30.842	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	24707m	3.000	ng/ul	0.03
60) Fluorene-d10	15.516	176	1067272	30.901	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.639	200	45309	6.394	ng/ul	0.00
73) Anthracene-d10	17.380	188	1737932	32.587	ng/ul	0.00
81) Pyrene-d10	19.616	212	1967770	34.138	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.639	264	1960974	33.135	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.334	88	104560	20.098	ng/uL#	91
52) Acenaphthene	14.586	153	44787	1.261	ng/ul	98

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

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