

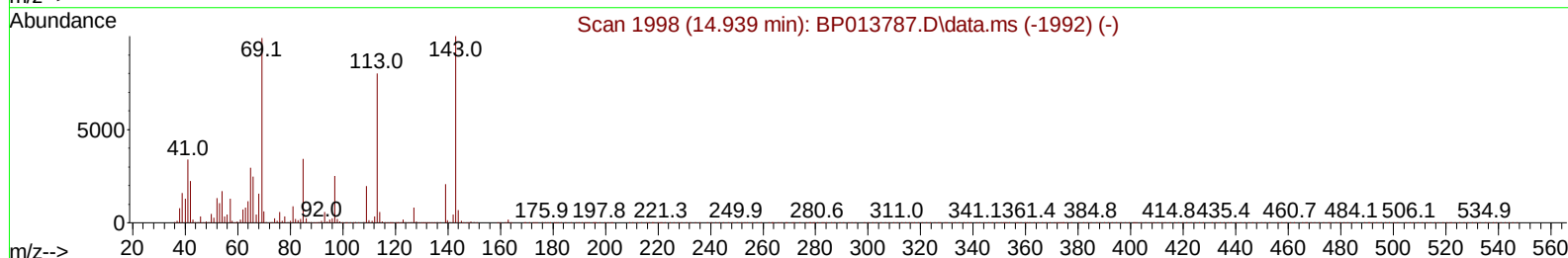
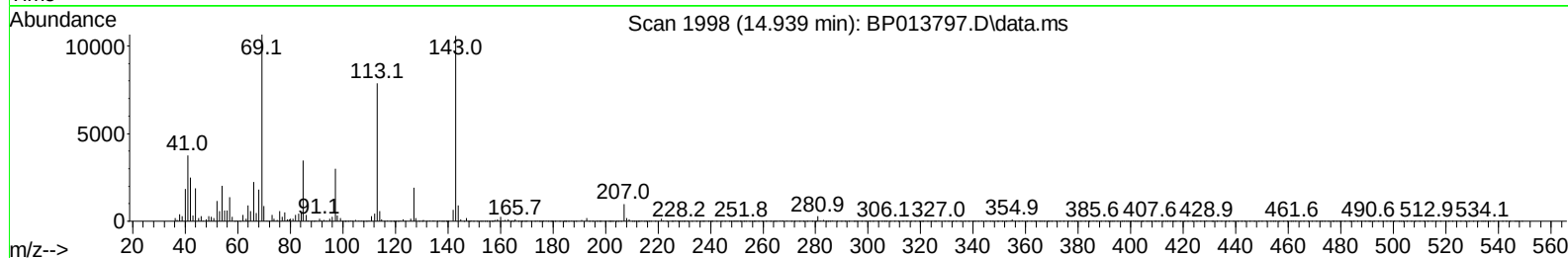
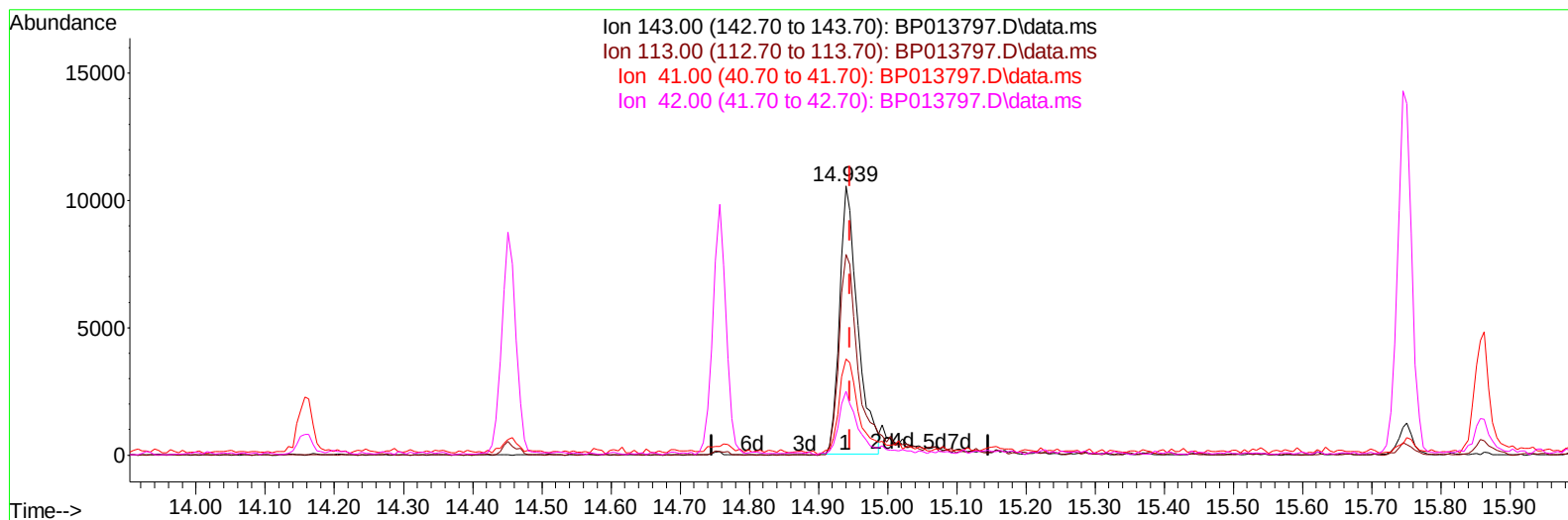
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020823\
 Data File : BP013797.D
 Acq On : 08 Feb 2023 15:14
 Operator : CG/JU
 Sample : 01422-04
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0B63

Manual IntegrationsAPPROVED

Quant Time: Feb 08 23:56:05 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020723.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Feb 08 00:37:14 2023
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/09/2023
 Supervised By :Sohil Jodhani 02/09/2023



TIC: BP013797.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.939min (-0.006) 2.77 ng/ul

response 19249

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	74.65
41.00	33.80	35.63
42.00	22.00	23.70

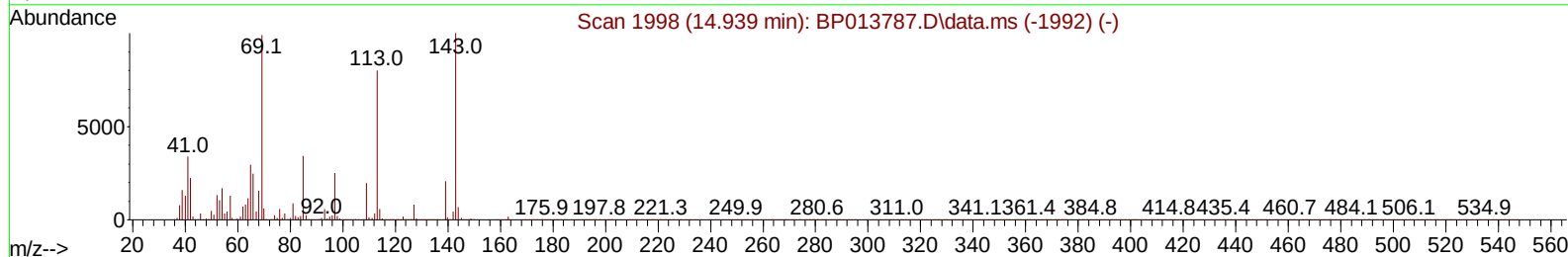
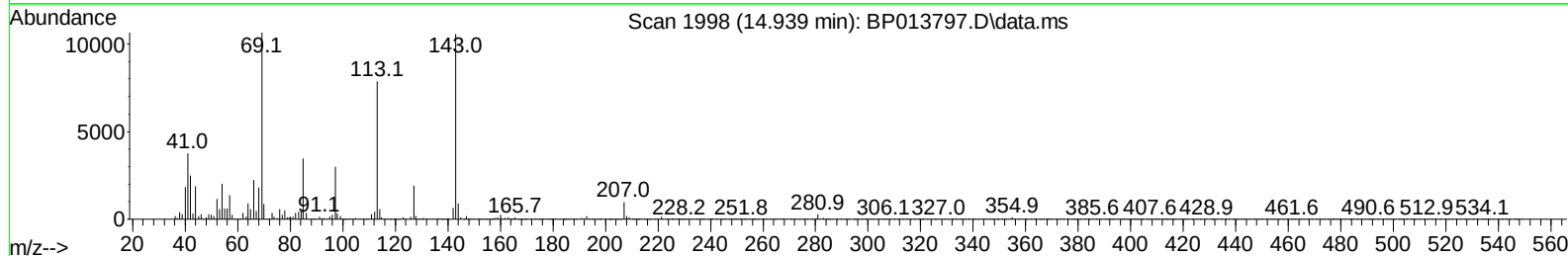
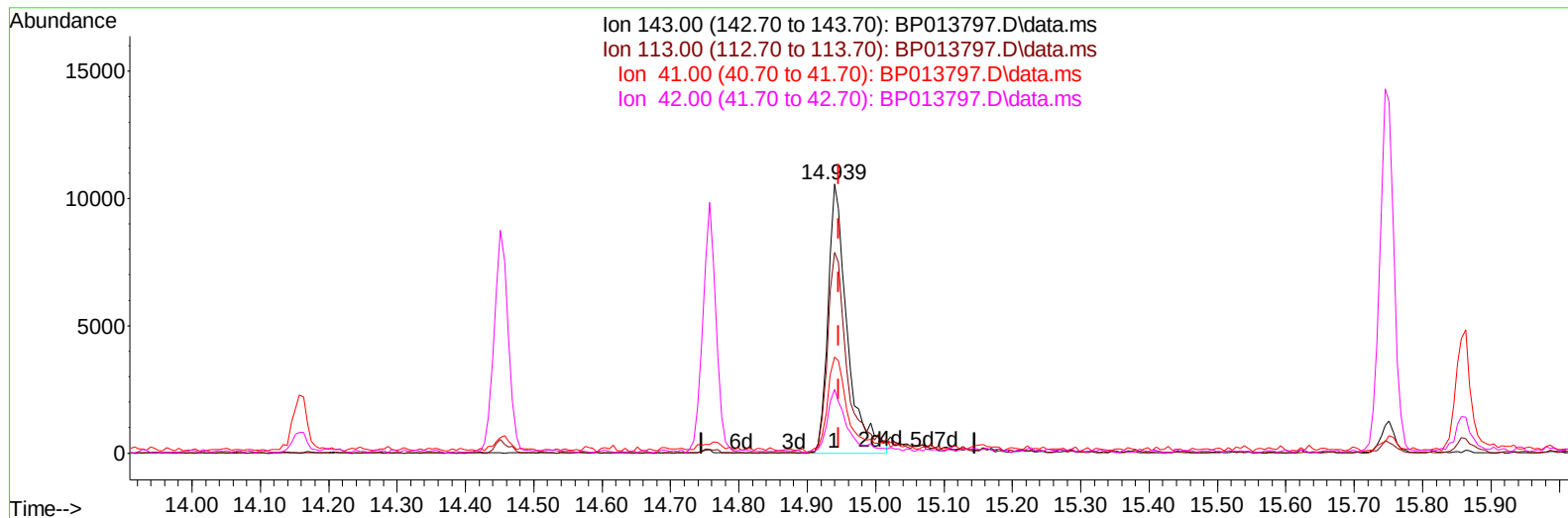
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(54) 4-Nitrophenol-d4 (S)

14.939min (-0.006) 2.97 ng/ul m

response 20614

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	78.80	74.65
41.00	33.80	35.63
42.00	22.00	23.70

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.122	152	158976	20.000	ng/ul	0.00
20) Naphthalene-d8	10.951	136	688214	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.757	164	495442	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.510	188	1196097	20.000	ng/ul	0.00
79) Chrysene-d12	21.586	240	1302998	20.000	ng/ul	0.00
88) Perylene-d12	24.151	264	1357486	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.469	96	16684	4.350	ng/uL	0.00
4) Pyridine-d5	3.911	84	55224	4.856	ng/ul	0.00
7) Phenol-d5	7.269	99	69256	4.889	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.440	67	215910	26.012	ng/ul	0.00
11) 2-Chlorophenol-d4	7.646	132	216705	20.783	ng/ul	0.00
15) 4-Methylphenol-d8	8.828	113	147434	12.735	ng/ul	0.00
21) Nitrobenzene-d5	9.293	128	133352	27.849	ng/ul	0.00
24) 2-Nitrophenol-d4	10.022	143	127834	23.122	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.563	165	266060	23.615	ng/ul	0.00
31) 4-Chloroaniline-d4	11.081	131	357608	22.392	ng/ul	0.00
46) Dimethylphthalate-d6	14.157	166	1129514	28.945	ng/ul	0.00
49) Acenaphthylene-d8	14.451	160	1176668	27.856	ng/ul	0.00
54) 4-Nitrophenol-d4	14.939	143	20614m	2.965	ng/ul	0.00
60) Fluorene-d10	15.751	176	979761	29.477	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.863	200	84240	11.480	ng/ul	0.00
73) Anthracene-d10	17.610	188	1866178	34.121	ng/ul	0.00
81) Pyrene-d10	19.827	212	2412172	32.645	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.986	264	2362528	34.733	ng/ul	0.00

Target Compounds Qvalue

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

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