

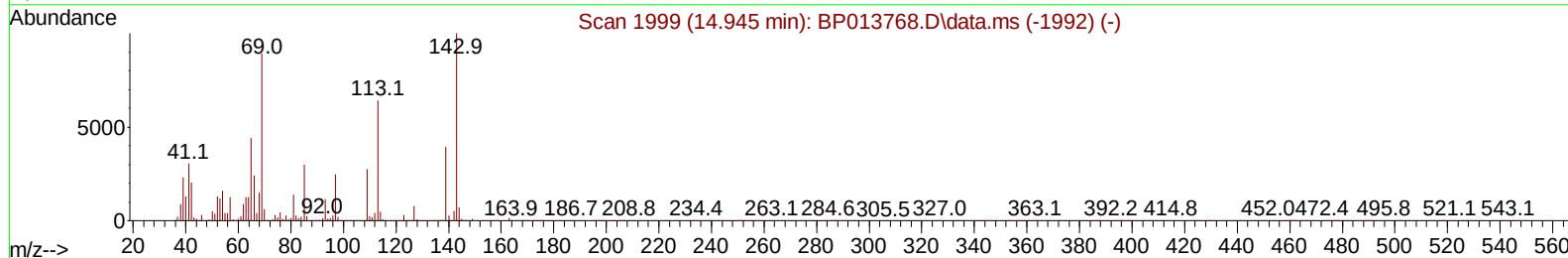
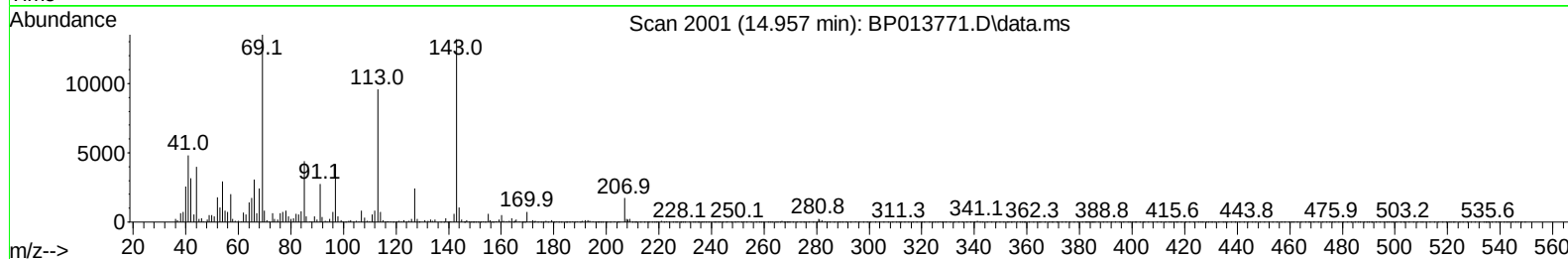
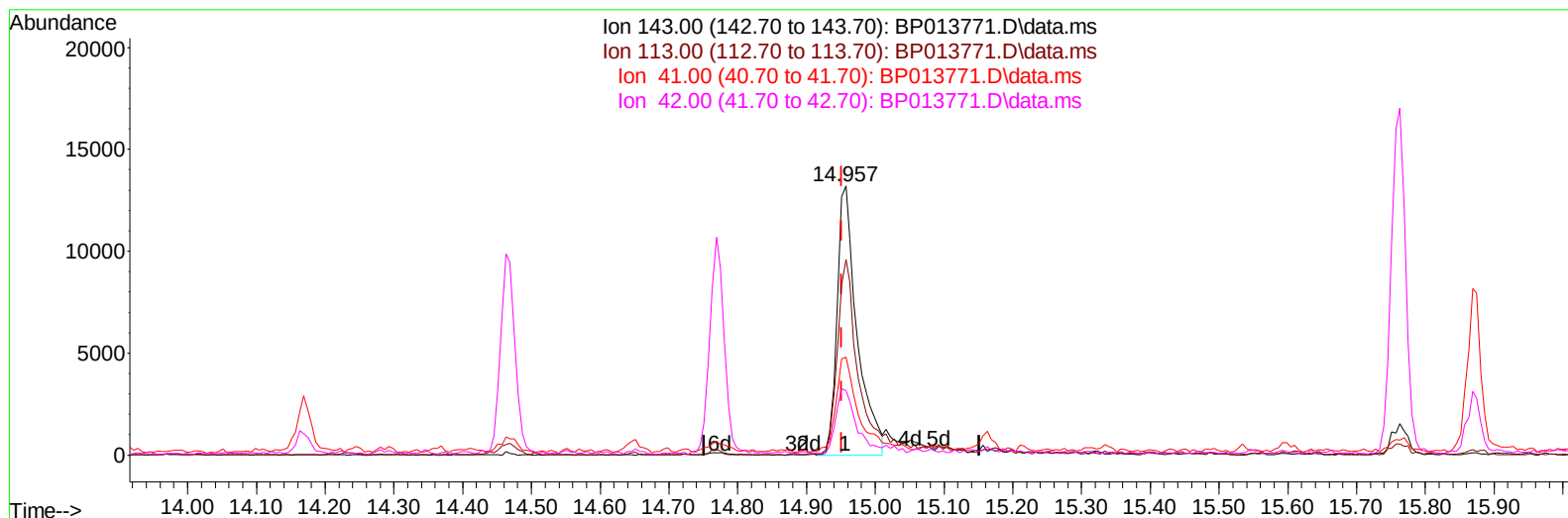
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020623\
 Data File : BP013771.D
 Acq On : 06 Feb 2023 13:58
 Operator : CG/JU
 Sample : 01381-22
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A44M5

Manual IntegrationsAPPROVED

Quant Time: Feb 06 23:54:35 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 03 23:44:43 2023
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/07/2023
 Supervised By :mohammad ahmed 02/07/2023



TIC: BP013771.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.957min (+ 0.006) 3.00 ng/ul

response 26758

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	69.30	72.72
41.00	33.60	36.55
42.00	21.80	23.94

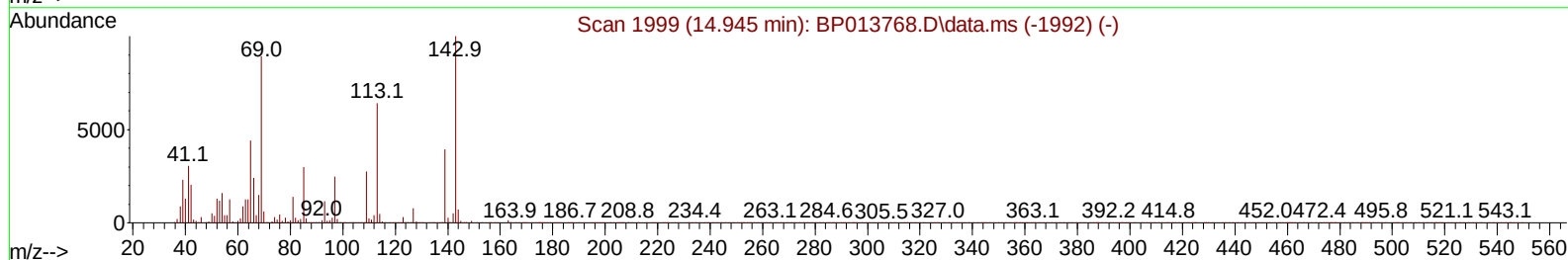
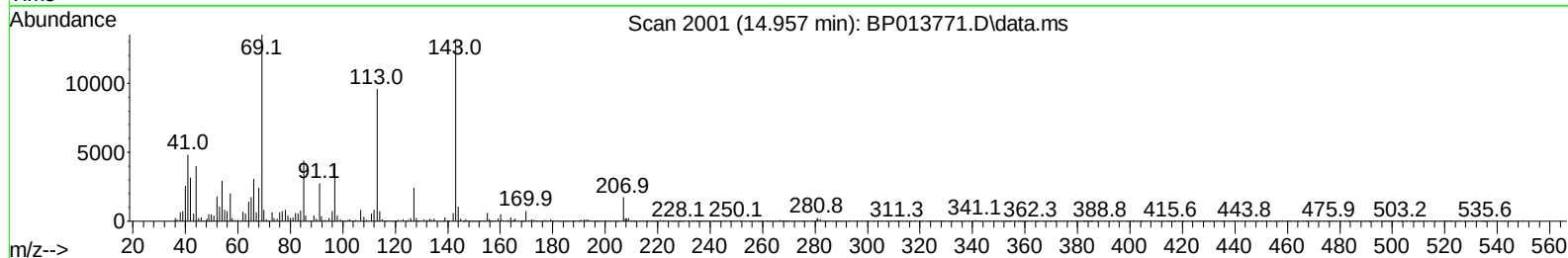
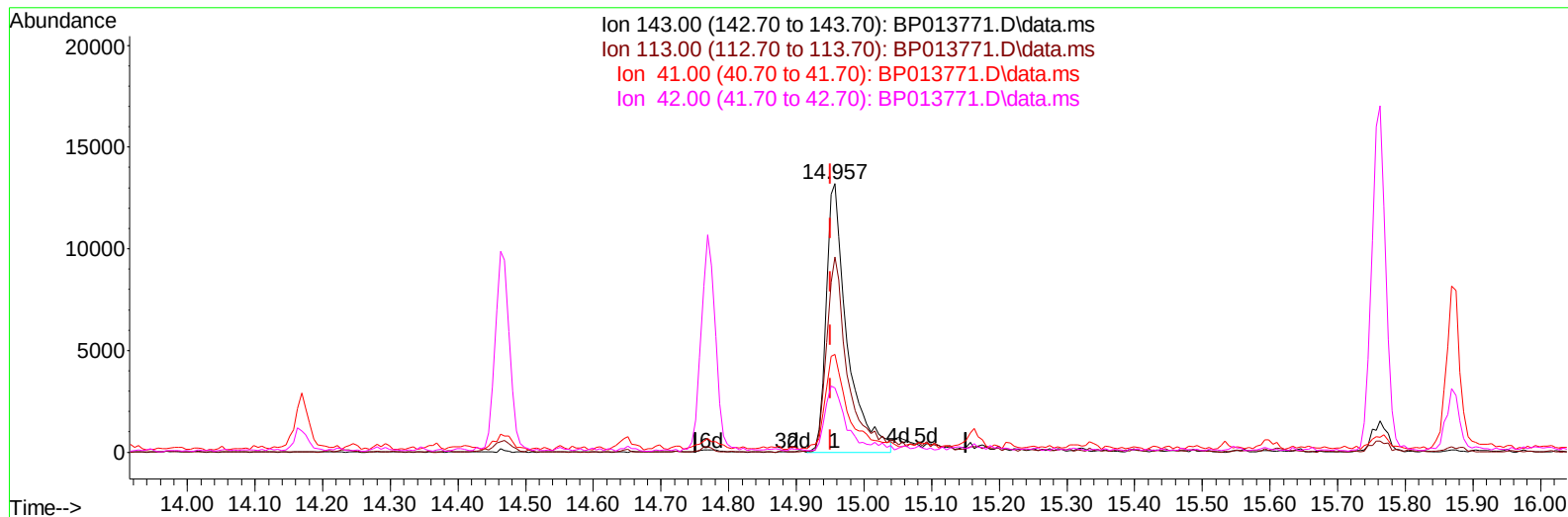
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TIC: BP013771.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.957min (+ 0.006) 3.17 ng/ul m

response 28209

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	69.30	72.72
41.00	33.60	36.55
42.00	21.80	23.94

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.134	152	202718	20.000	ng/ul	0.00
20) Naphthalene-d8	10.963	136	907103	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.769	164	670505	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.522	188	1707620	20.000	ng/ul	0.00
79) Chrysene-d12	21.598	240	1815519	20.000	ng/ul	0.00
88) Perylene-d12	24.174	264	1815933	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.469	96	21801	4.560	ng/uL	0.00
4) Pyridine-d5	3.916	84	77599	5.656	ng/ul	0.00
7) Phenol-d5	7.281	99	101879	5.718	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.452	67	292585	28.000	ng/ul	0.00
11) 2-Chlorophenol-d4	7.657	132	285101	21.829	ng/ul	0.00
15) 4-Methylphenol-d8	8.840	113	203935	14.358	ng/ul	0.00
21) Nitrobenzene-d5	9.310	128	165914	28.672	ng/ul	0.00
24) 2-Nitrophenol-d4	10.034	143	183766	27.250	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.575	165	387863	24.595	ng/ul	0.00
31) 4-Chloroaniline-d4	11.098	131	259057	12.438	ng/ul	0.00
46) Dimethylphthalate-d6	14.169	166	1521884	28.794	ng/ul	0.00
49) Acenaphthylene-d8	14.469	160	1597790	27.362	ng/ul	0.00
54) 4-Nitrophenol-d4	14.957	143	28209m	3.166	ng/ul	0.00
60) Fluorene-d10	15.763	176	1429379	31.376	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.875	200	189781	19.832	ng/ul	0.00
73) Anthracene-d10	17.622	188	2722063	34.198	ng/ul	0.00
81) Pyrene-d10	19.839	212	3572740	35.283	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.010	264	3224212	34.781	ng/ul	0.00

Target Compounds Qvalue

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

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