

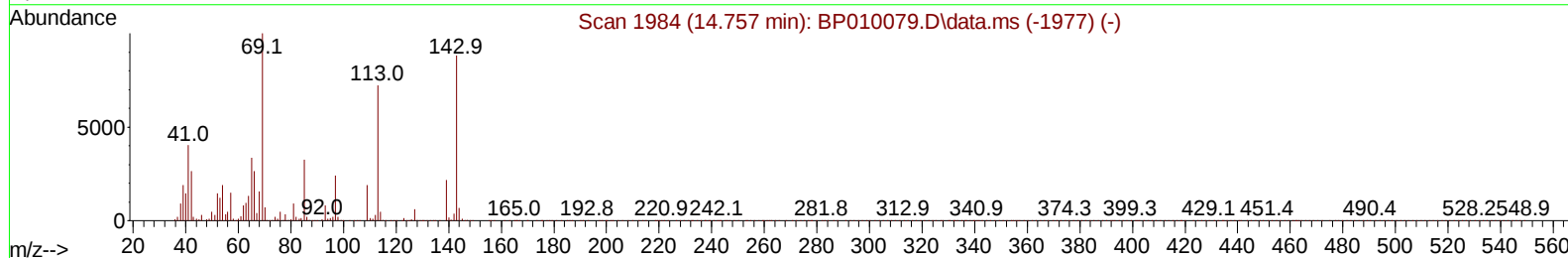
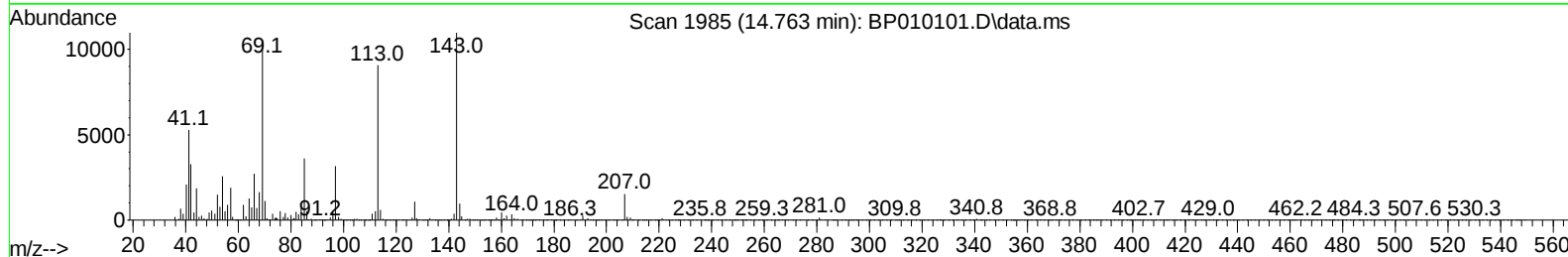
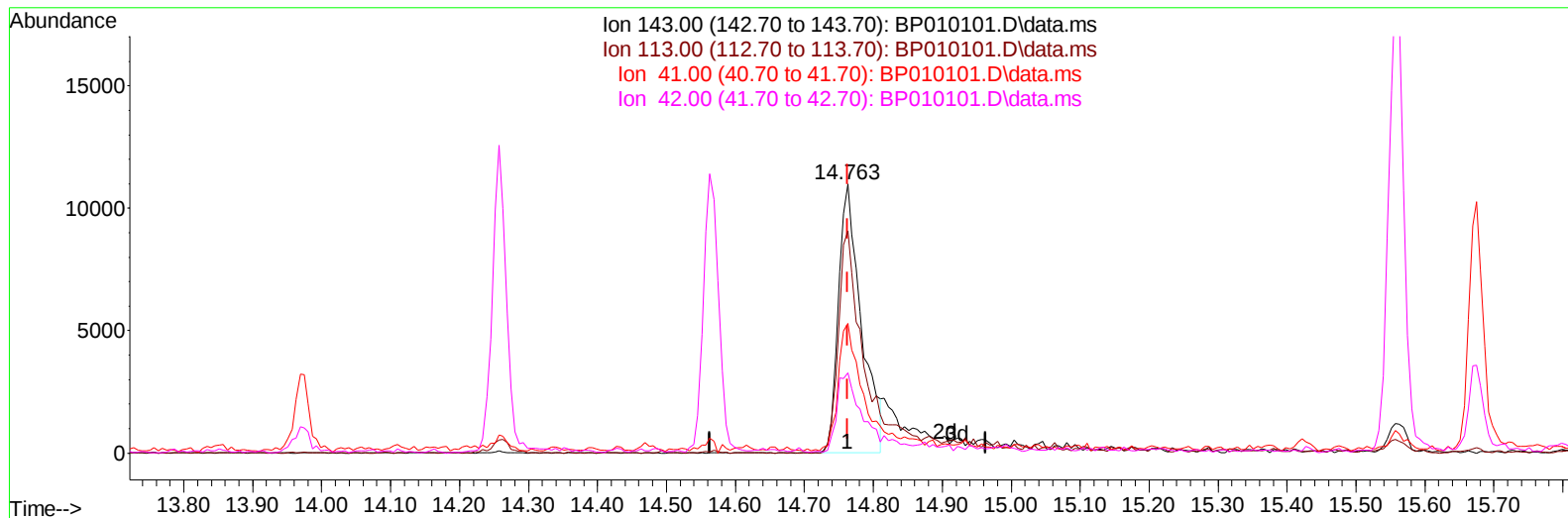
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
Data File : BP010101.D
Acq On : 26 Apr 2022 01:31
Operator : CG/JU
Sample : N2475-13
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A0BN0

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/26/2022
Supervised By :Yogesh Patel 04/28/2022

Quant Time: Apr 26 02:51:56 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 14:49:38 2022
Response via : Initial Calibration



TIC: BP010101.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.000) 3.31 ng/ul

response 25184

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	82.62#
41.00	23.20	48.24#
42.00	18.00	29.72#

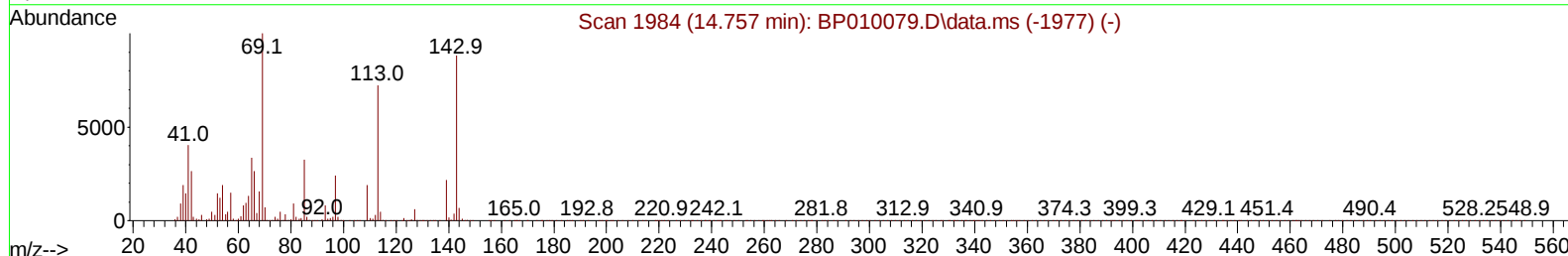
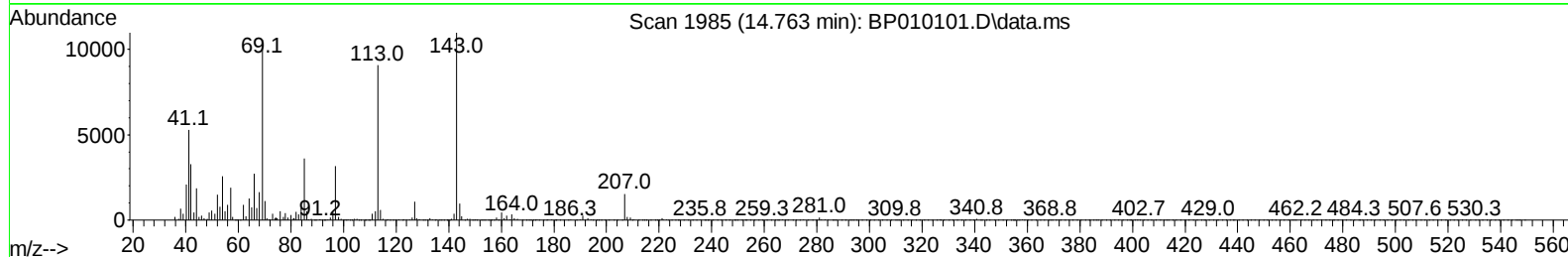
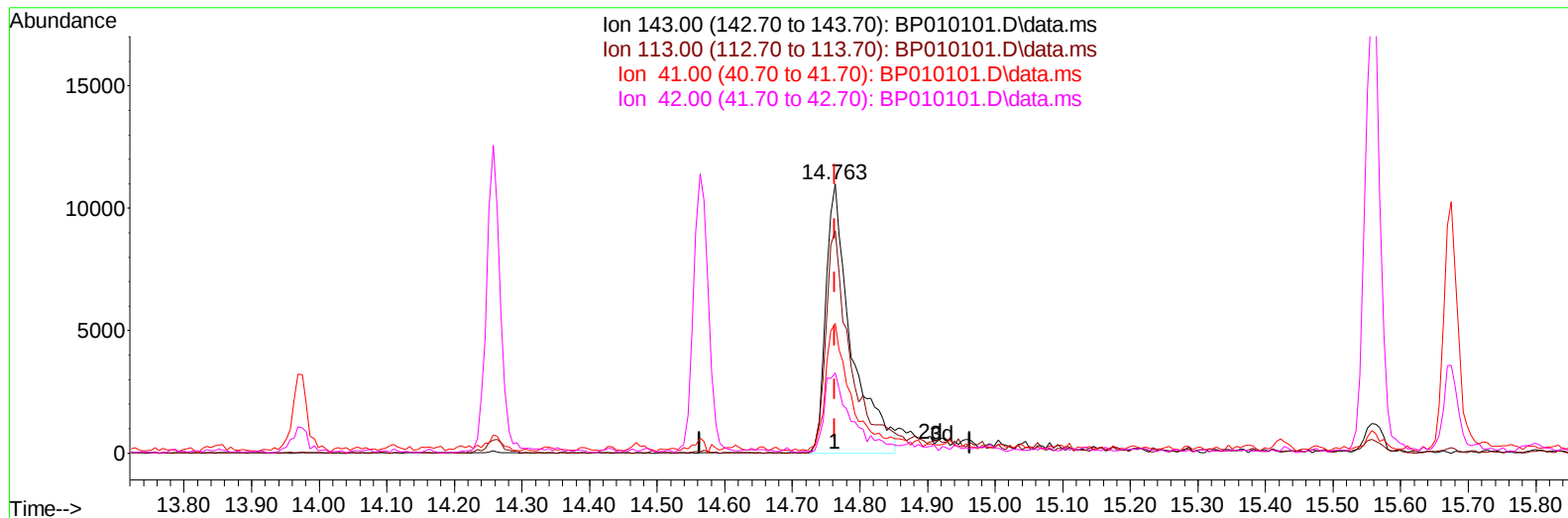
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TIC: BP010101.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.000) 3.79 ng/ul m

response 28839

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	82.62#
41.00	23.20	48.24#
42.00	18.00	29.72#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	180867	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	780299	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	531659	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.316	188	1091503	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.398	240	880283	20.000	ng/ul	#-0.01
88) Perylene-d12	23.851	264	666571	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	18949	4.569	ng/uL	0.00
4) Pyridine-d5	3.787	84	72643	6.226	ng/ul	0.00
7) Phenol-d5	7.093	99	93406	5.678	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.258	67	286570	29.030	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.463	132	269686	22.135	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	189780	13.923	ng/ul	-0.01
21) Nitrobenzene-d5	9.093	128	180831	29.138	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	190225	28.066	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	311642	23.618	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	426035	22.519	ng/ul	-0.01
46) Dimethylphthalate-d6	13.975	166	1316054	31.822	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1418562	28.341	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	28839m	3.788	ng/ul	0.00
60) Fluorene-d10	15.557	176	1099794	30.914	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	162576	23.556	ng/ul	0.00
73) Anthracene-d10	17.416	188	1781625	33.739	ng/ul	0.00
81) Pyrene-d10	19.645	212	2024770	40.445	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.692	264	1219995	34.911	ng/ul	0.00

Target Compounds Qvalue

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

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