

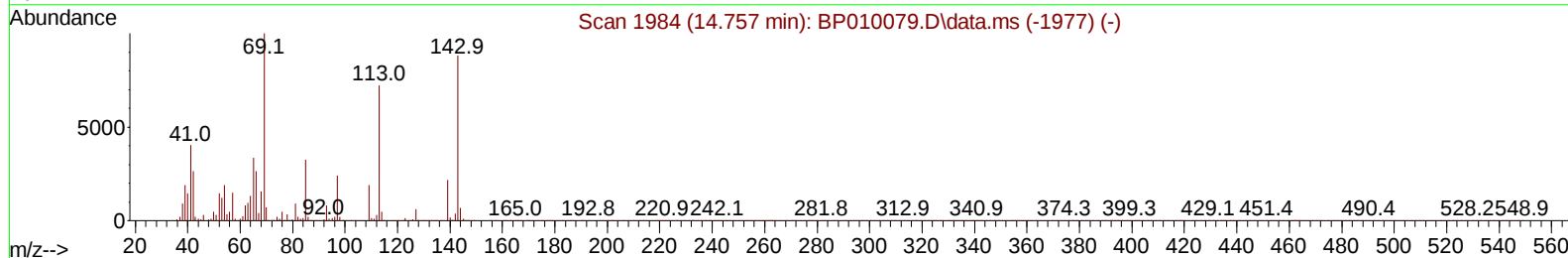
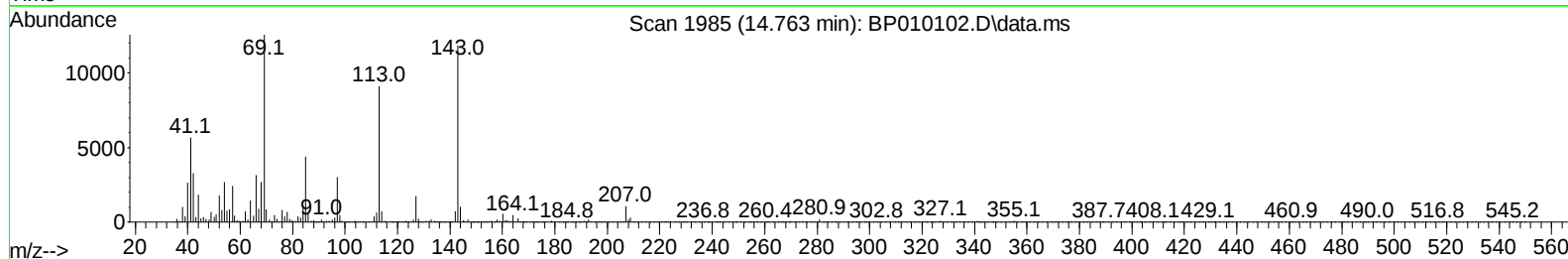
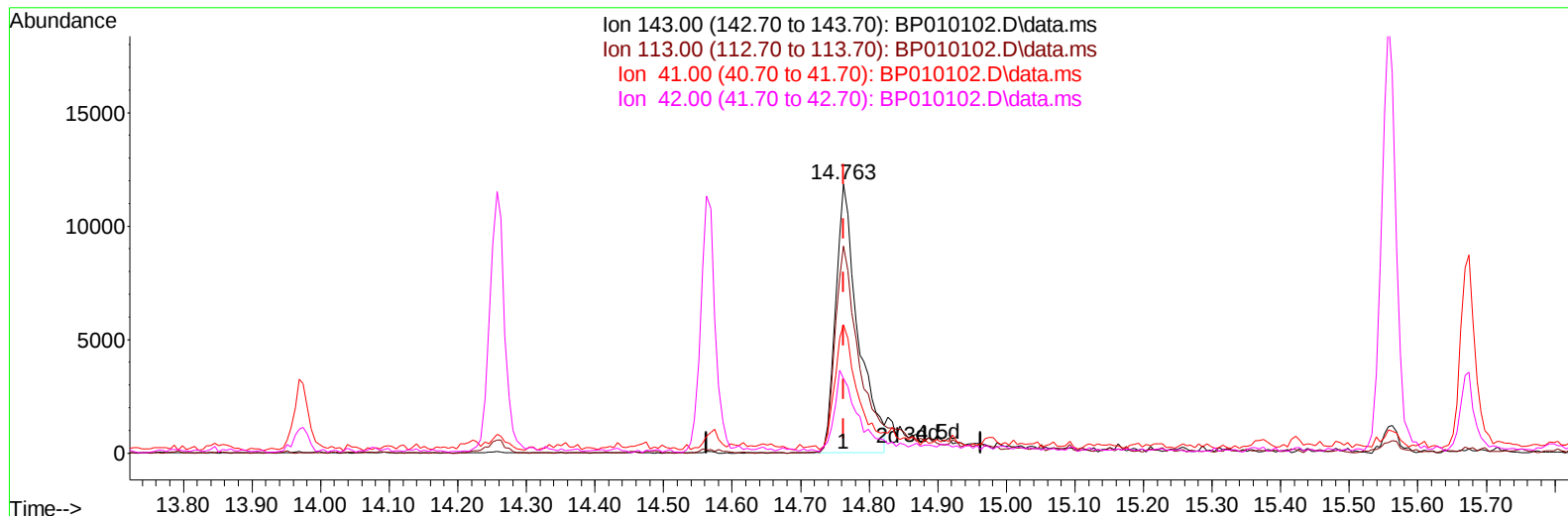
Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042522\
 Data File : BP010102.D
 Acq On : 26 Apr 2022 02:04
 Operator : CG/JU
 Sample : N2475-14
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A0BN4

Manual IntegrationsAPPROVED

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

Quant Time: Apr 26 02:52:11 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: BP010102.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.000) 3.72 ng/u1

response 26802

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	76.90#
41.00	23.20	47.77#
42.00	18.00	27.81#

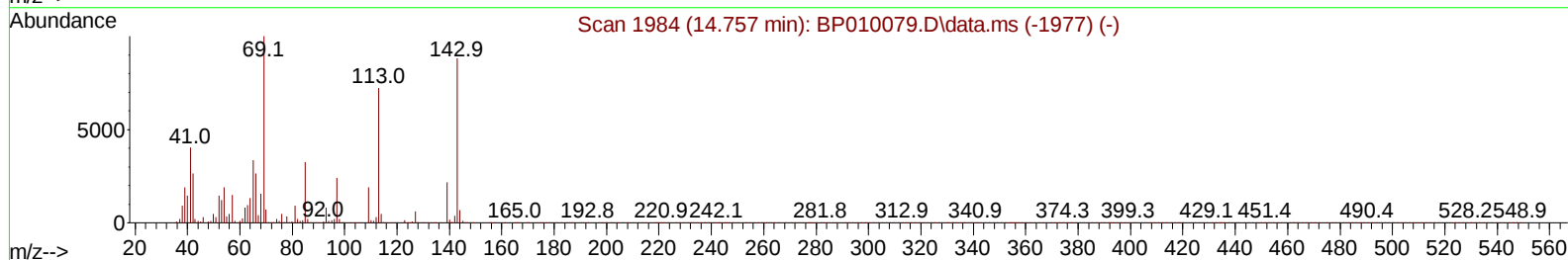
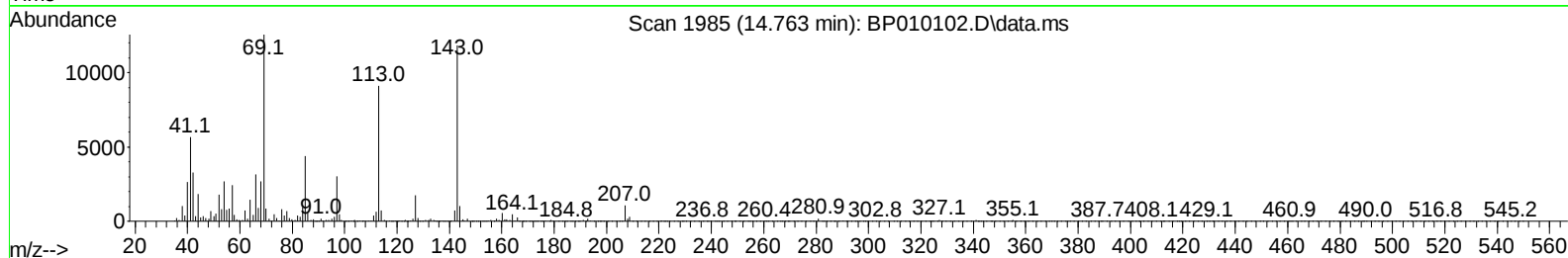
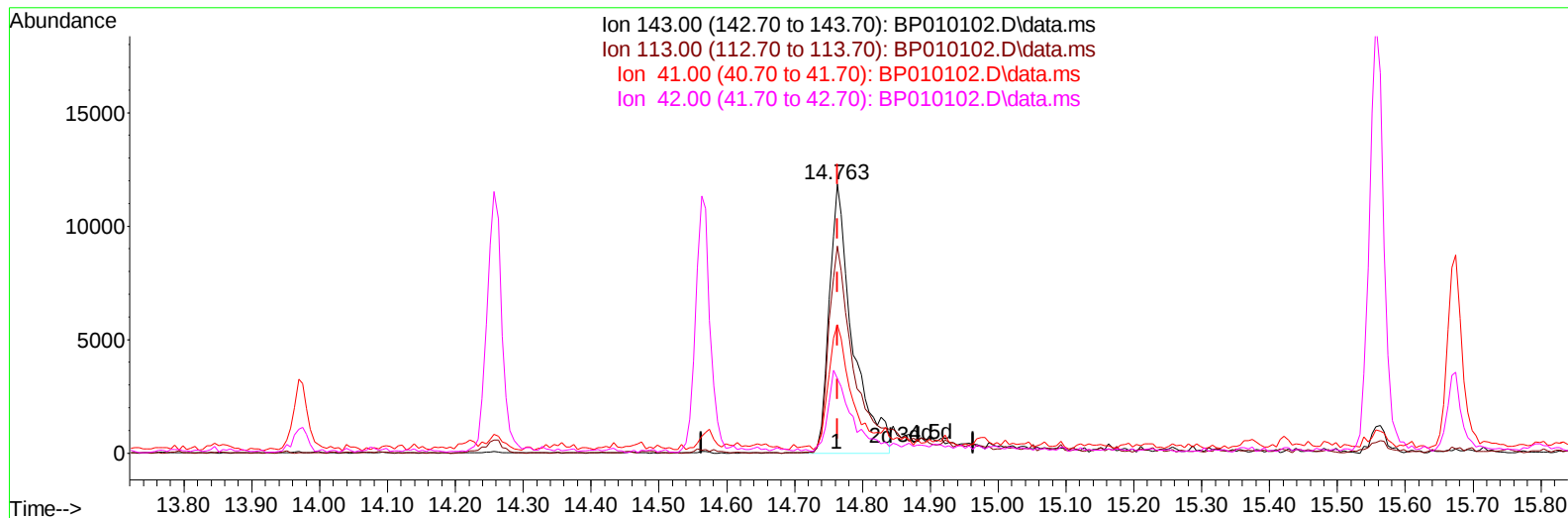
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TIC: BP010102.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (-0.000) 3.91 ng/ul m

response 28167

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	132.70	76.90#
41.00	23.20	47.77#
42.00	18.00	27.81#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.934	152	182779	20.000	ng/ul	0.00
20) Naphthalene-d8	10.740	136	766001	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.569	164	503518	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.316	188	970783	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.398	240	647293	20.000	ng/ul	#-0.01
88) Perylene-d12	23.851	264	591319	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	21707	5.179	ng/uL	0.00
4) Pyridine-d5	3.787	84	85729	7.270	ng/ul	0.00
7) Phenol-d5	7.093	99	104105	6.262	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.257	67	292231	29.293	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.463	132	282475	22.942	ng/ul	0.00
15) 4-Methylphenol-d8	8.640	113	202542	14.704	ng/ul	-0.01
21) Nitrobenzene-d5	9.093	128	185562	30.458	ng/ul	0.00
24) 2-Nitrophenol-d4	9.816	143	194639	29.253	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.357	165	328353	25.349	ng/ul	0.00
31) 4-Chloroaniline-d4	10.869	131	399823	21.528	ng/ul	-0.01
46) Dimethylphthalate-d6	13.975	166	1252642	31.981	ng/ul	0.00
49) Acenaphthylene-d8	14.257	160	1402034	29.576	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	28167m	3.906	ng/ul	0.00
60) Fluorene-d10	15.557	176	1055883	31.339	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.675	200	145701	23.736	ng/ul	0.00
73) Anthracene-d10	17.416	188	1564096	33.303	ng/ul	0.00
81) Pyrene-d10	19.645	212	1658483	45.053	ng/ul	-0.01
92) Benzo(a)pyrene-d12	23.692	264	1107745	35.733	ng/ul	0.00

Target Compounds Qvalue

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

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