

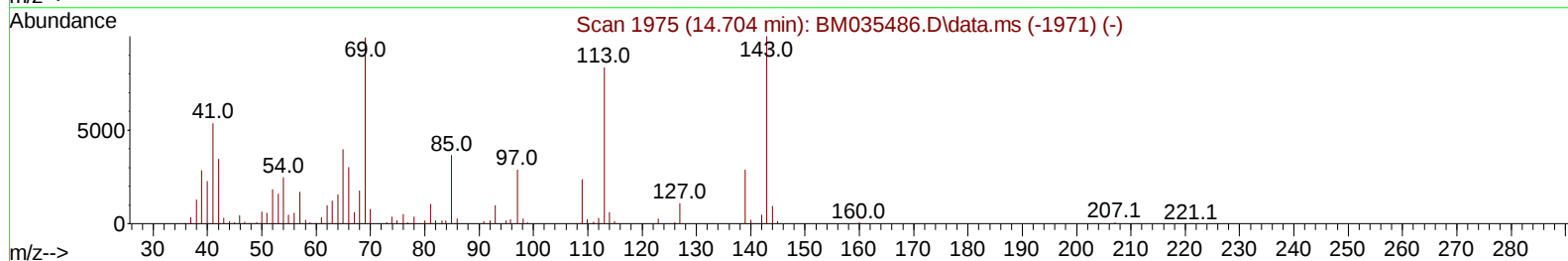
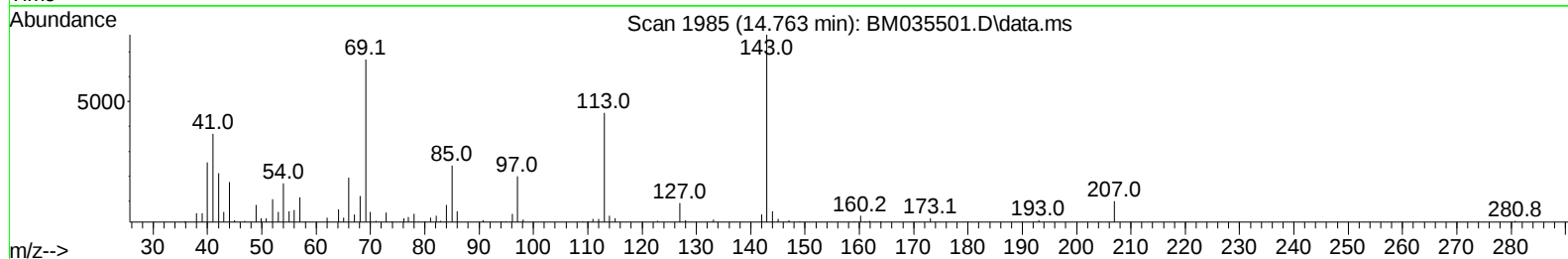
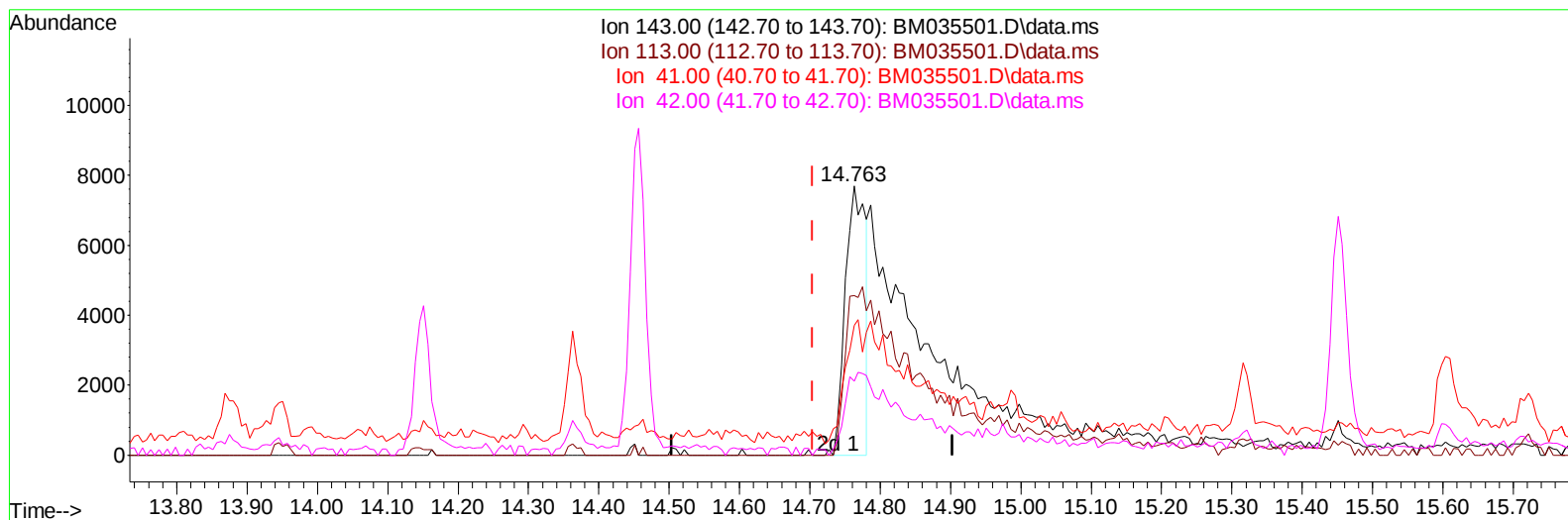
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061722\
Data File : BM035501.D
Acq On : 17 Jun 2022 18:54
Operator : CG/JU
Sample : N3226-18
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EW4N8

Manual IntegrationsAPPROVED

Quant Time: Jun 17 23:02:49 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM061622.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Jun 17 01:58:49 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/18/2022
Supervised By :Yogesh Patel 06/22/2022



TIC: BM035501.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.059) 2.99 ng/u1

response 15321

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	59.26
41.00	41.00	48.26
42.00	29.30	27.43

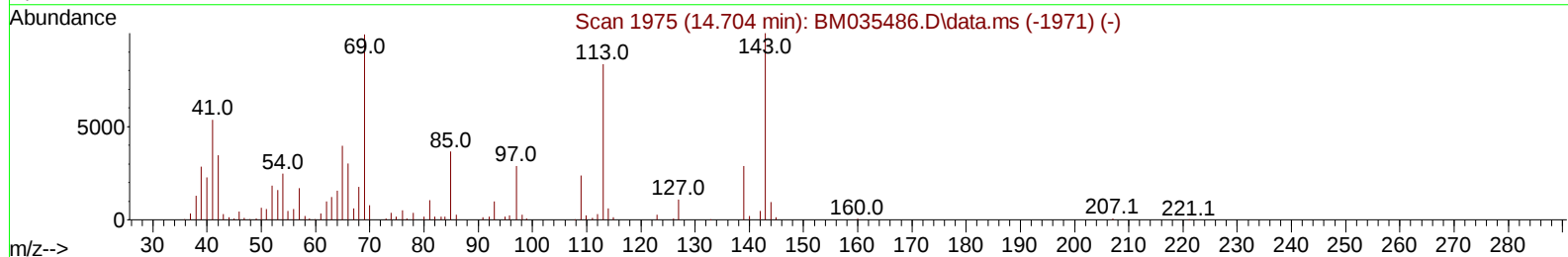
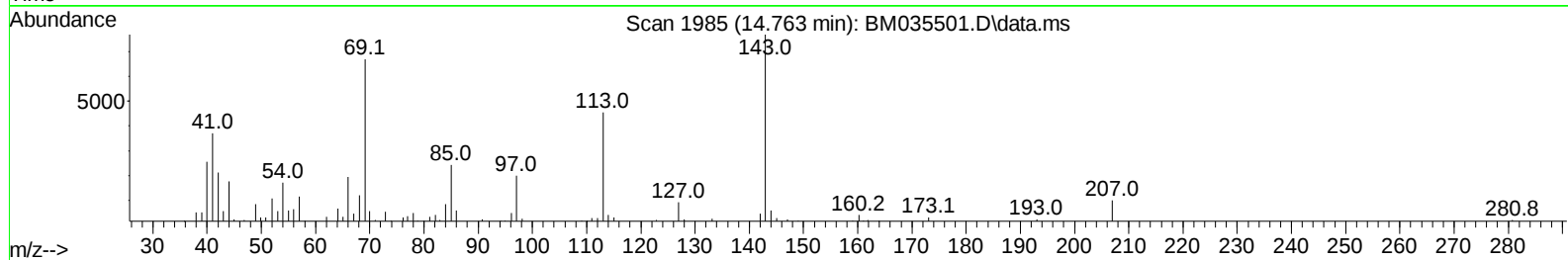
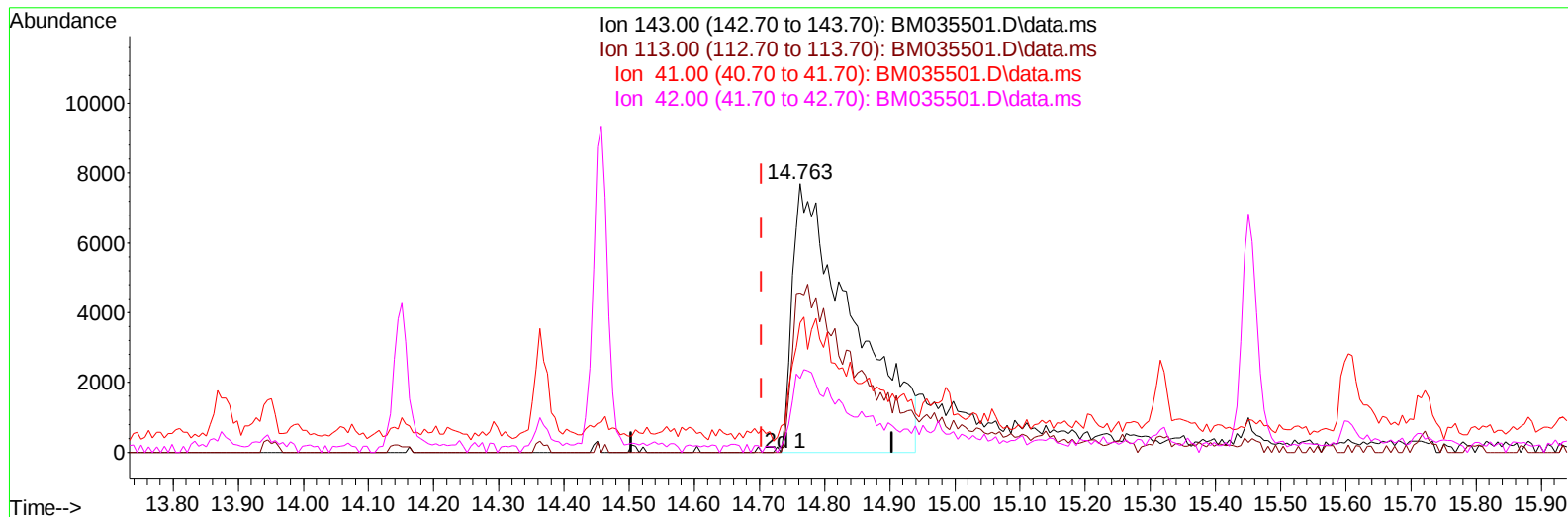
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(54) 4-Nitrophenol-d4 (S)

14.763min (+ 0.059) 9.49 ng/ul m

response 48703

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	59.26
41.00	41.00	48.26
42.00	29.30	27.43

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	194162	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	822517	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	491042	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1014239	20.000	ng/ul	0.00
79) Chrysene-d12	21.392	240	1022253	20.000	ng/ul	0.00
88) Perylene-d12	23.733	264	1096680	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	13346	2.744	ng/uL	0.00
4) Pyridine-d5	3.699	84	111026	9.268	ng/ul	0.01
7) Phenol-d5	7.010	99	190622	12.810	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.151	67	127522	13.781	ng/ul	0.00
11) 2-Chlorophenol-d4	7.357	132	185352	14.381	ng/ul	0.00
15) 4-Methylphenol-d8	8.551	113	158057	12.688	ng/ul	0.01
21) Nitrobenzene-d5	8.981	128	81121	13.249	ng/ul	0.00
24) 2-Nitrophenol-d4	9.710	143	85119	12.893	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.269	165	134287	11.980	ng/ul	0.02
31) 4-Chloroaniline-d4	10.769	131	212666	12.058	ng/ul	0.01
46) Dimethylphthalate-d6	13.875	166	516818	15.473	ng/ul	0.00
49) Acenaphthylene-d8	14.151	160	601725	14.537	ng/ul	0.00
54) 4-Nitrophenol-d4	14.763	143	48703m	9.492	ng/ul	0.06
60) Fluorene-d10	15.451	176	453508	15.681	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	48473	8.016	ng/ul	0.01
73) Anthracene-d10	17.304	188	738078	16.190	ng/ul	0.00
81) Pyrene-d10	19.598	212	886400	17.637	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.580	264	938233	17.738	ng/ul	0.00

Target Compounds Qvalue

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

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