

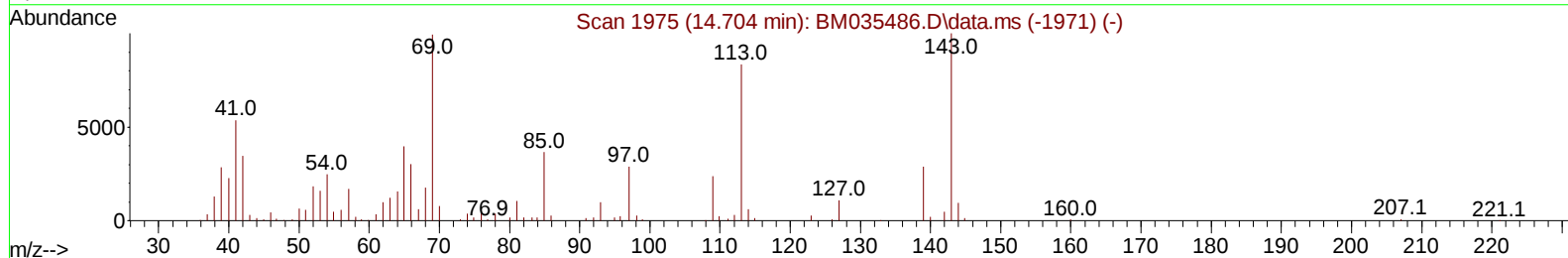
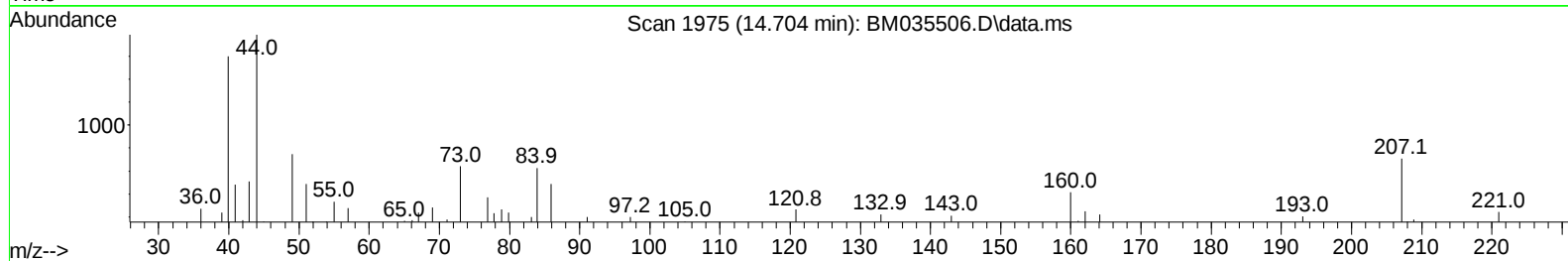
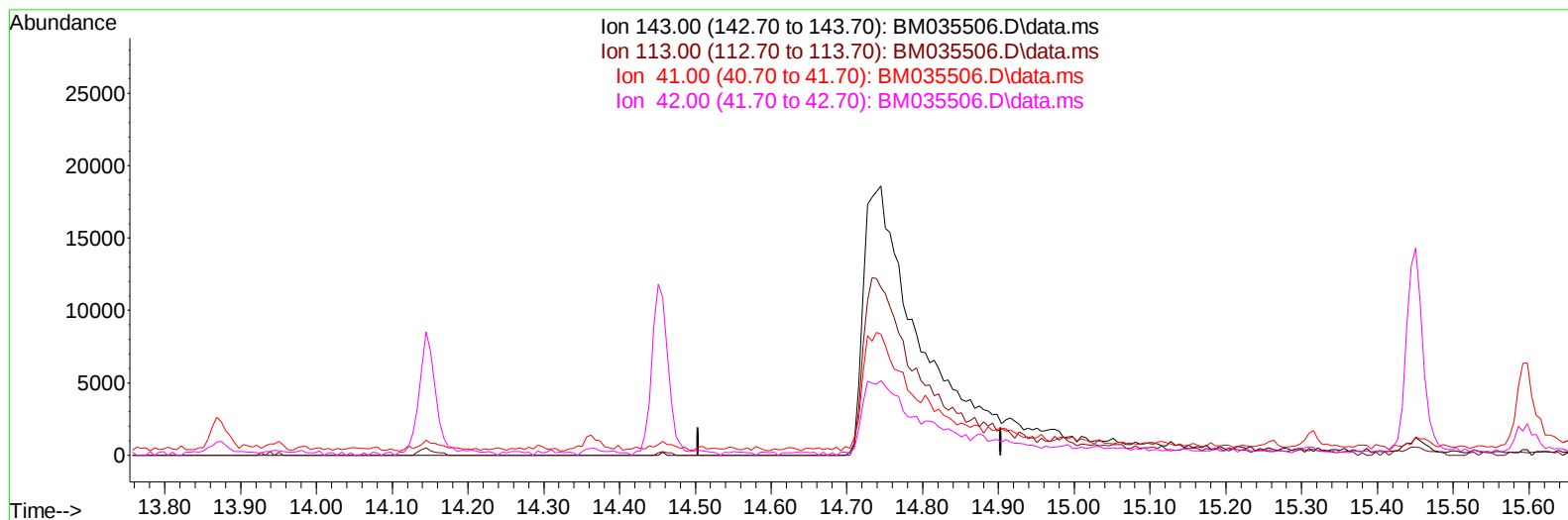
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061722\
 Data File : BM035506.D
 Acq On : 17 Jun 2022 22:35
 Operator : CG/JU
 Sample : N3226-12
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4N2

Manual IntegrationsAPPROVED

Quant Time: Jun 18 00:12:37 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: BM035506.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.704min (-14.704) 0.00 ng/u1

response 0

Ion	Exp%	Act%
143.00	100.00	0.00
113.00	61.40	0.00#
41.00	41.00	0.00#
42.00	29.30	0.00#

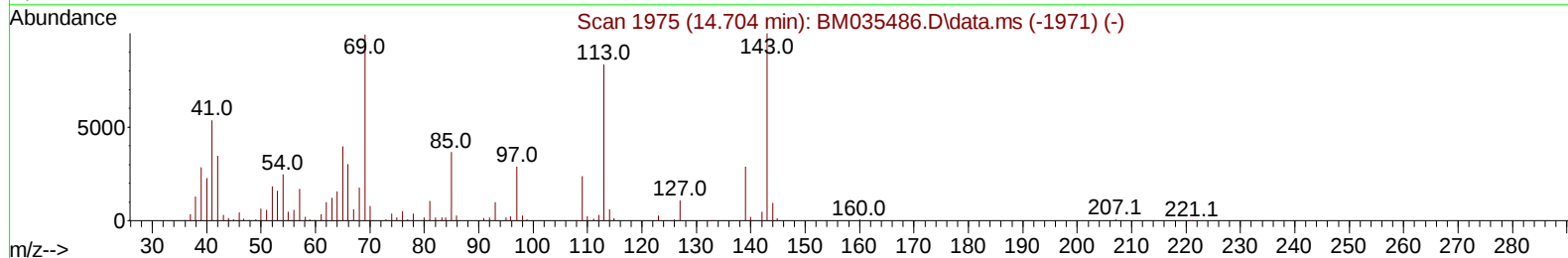
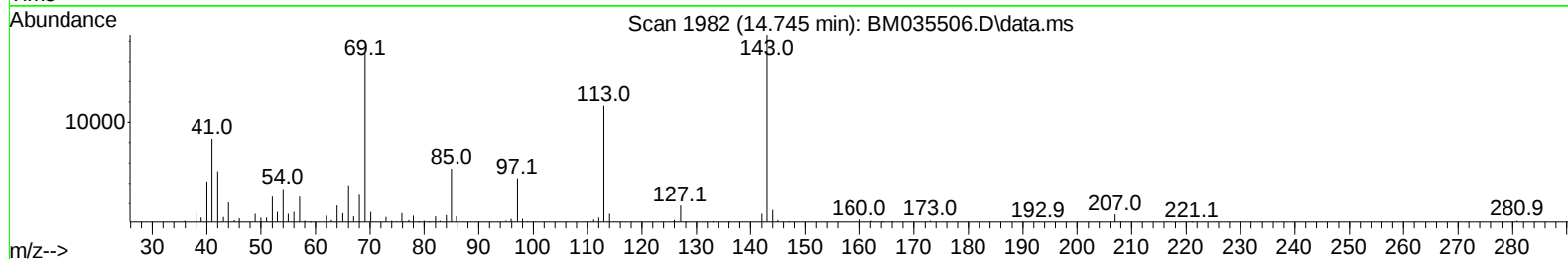
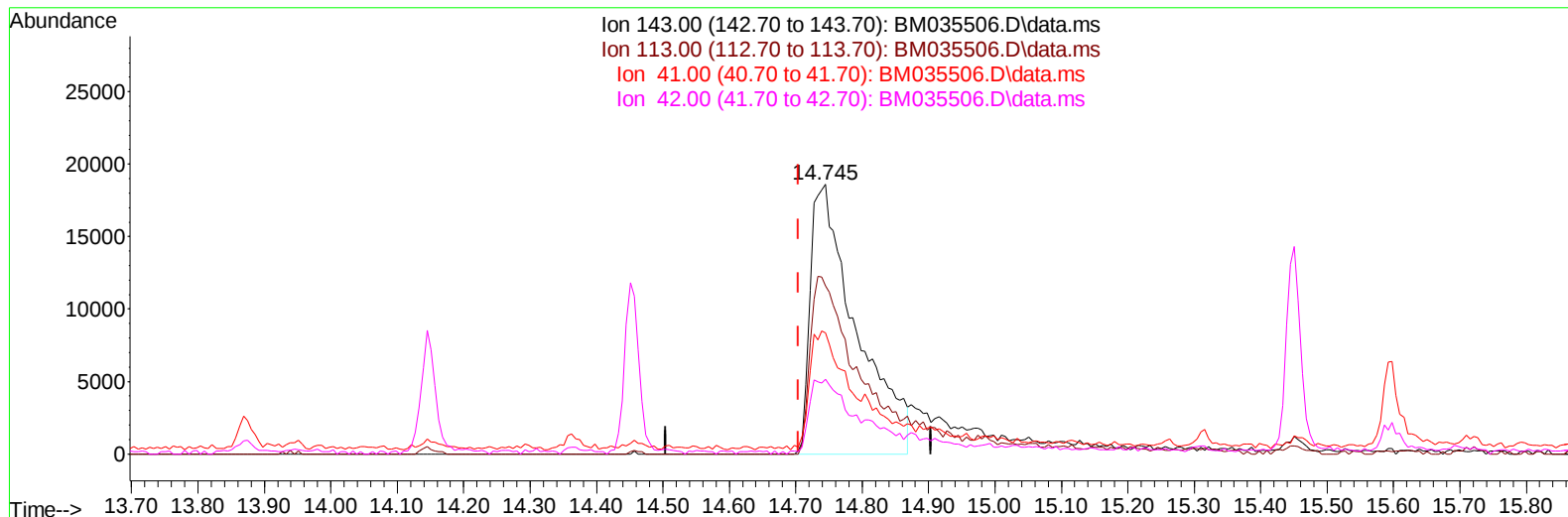
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TIC: BM035506.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.745min (+ 0.041) 14.93 ng/ul m

response 88931

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	62.48
41.00	41.00	44.84
42.00	29.30	27.64

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 Sample : N3226-12
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.810	152	219122	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	945338	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	570064	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1156856	20.000	ng/ul	0.00
79) Chrysene-d12	21.386	240	988862	20.000	ng/ul	0.00
88) Perylene-d12	23.727	264	1032471	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.269	96	20608	3.754	ng/uL	0.00
4) Pyridine-d5	3.693	84	221753	16.402	ng/ul	0.00
7) Phenol-d5	7.004	99	329008	19.591	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.145	67	229072	21.935	ng/ul	0.00
11) 2-Chlorophenol-d4	7.351	132	315975	21.723	ng/ul	0.00
15) 4-Methylphenol-d8	8.545	113	259102	18.431	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	144376	20.517	ng/ul	0.00
24) 2-Nitrophenol-d4	9.704	143	160931	21.209	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	260805	20.244	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	379880	18.740	ng/ul	0.00
46) Dimethylphthalate-d6	13.874	166	918325	23.682	ng/ul	0.00
49) Acenaphthylene-d8	14.145	160	1103623	22.967	ng/ul	0.00
54) 4-Nitrophenol-d4	14.745	143	88931m	14.930	ng/ul	0.04
60) Fluorene-d10	15.451	176	811846	24.180	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	102253	14.826	ng/ul	0.00
73) Anthracene-d10	17.298	188	1164834	22.401	ng/ul	0.00
81) Pyrene-d10	19.592	212	1293362	26.604	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.574	264	1285923	25.823	ng/ul	0.00

Target Compounds Qvalue

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

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