

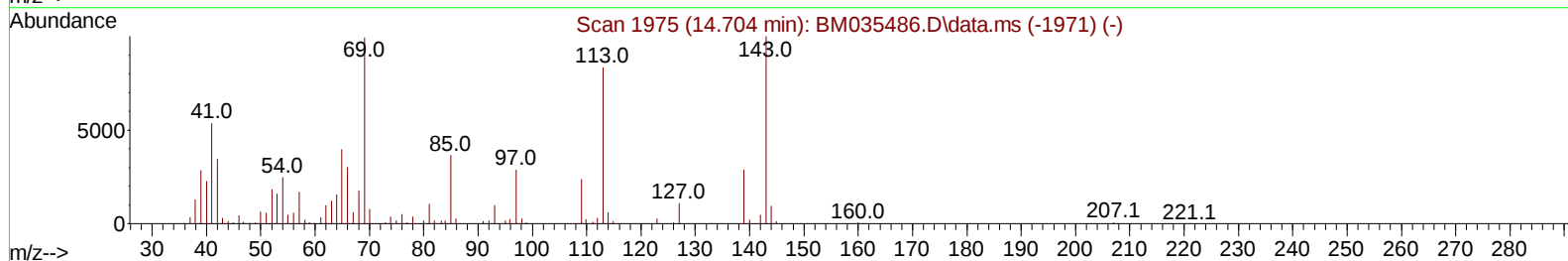
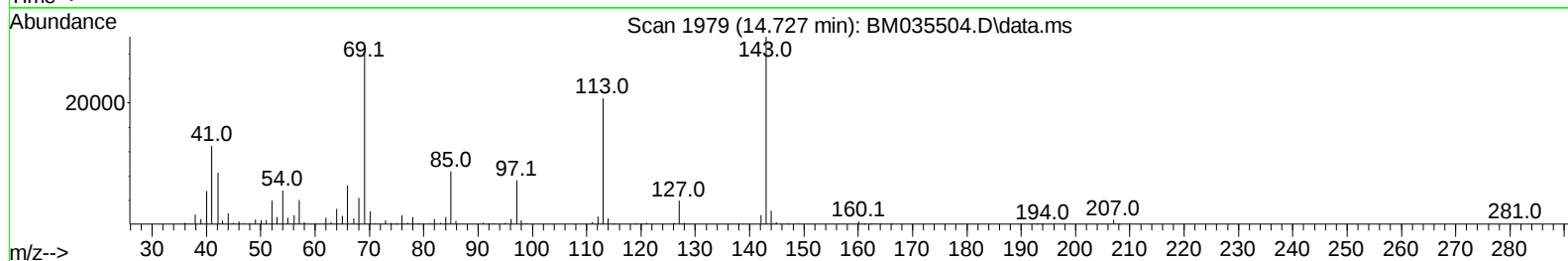
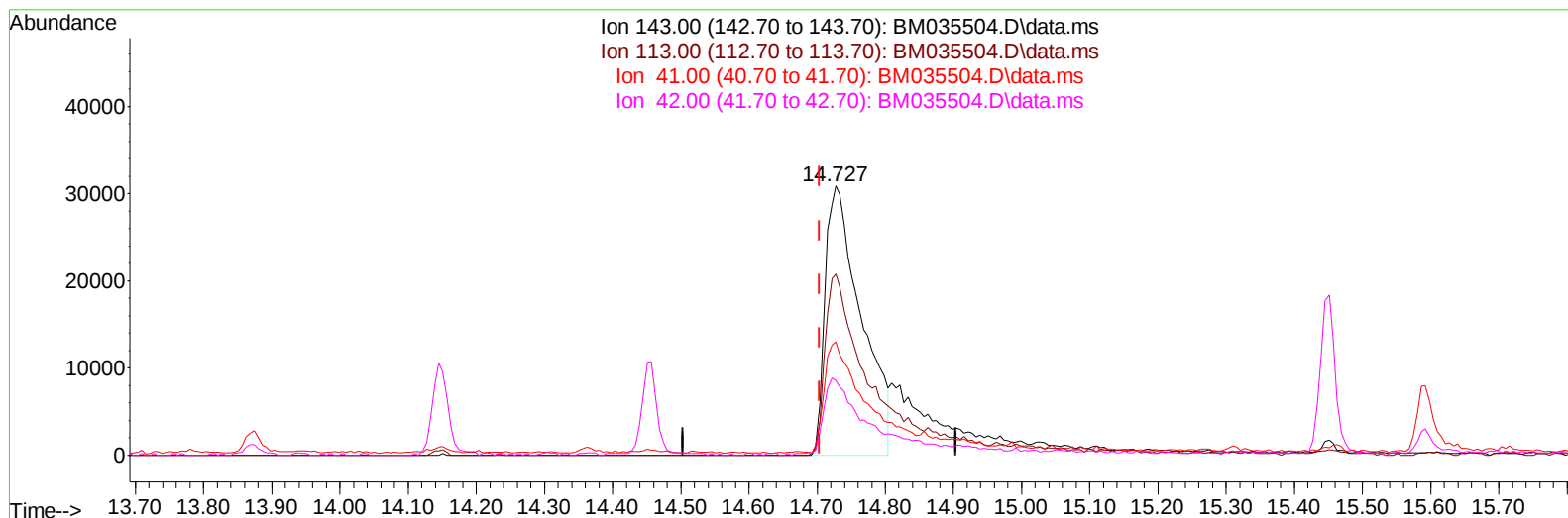
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061722\
 Data File : BM035504.D
 Acq On : 17 Jun 2022 21:22
 Operator : CG/JU
 Sample : N3226-05
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 EW4L6

Manual IntegrationsAPPROVED

Quant Time: Jun 17 23:03:17 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: BM035504.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.727min (+ 0.024) 19.15 ng/u1

response 113287

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	67.31
41.00	41.00	42.11
42.00	29.30	27.74

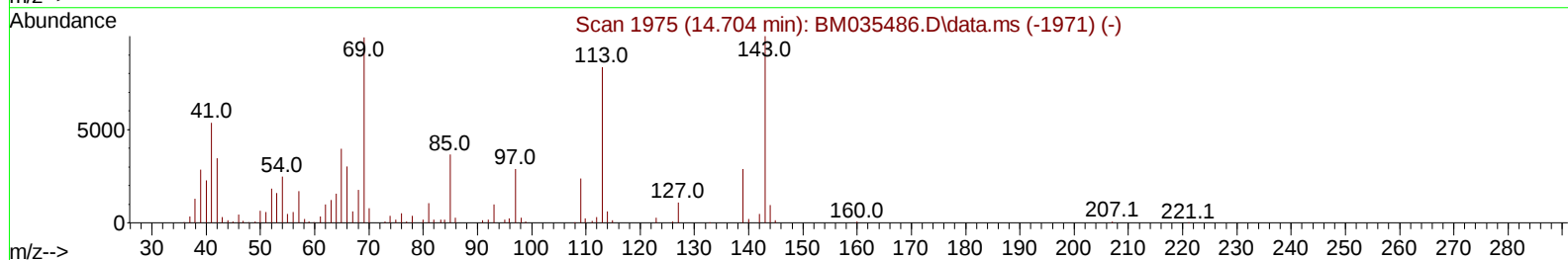
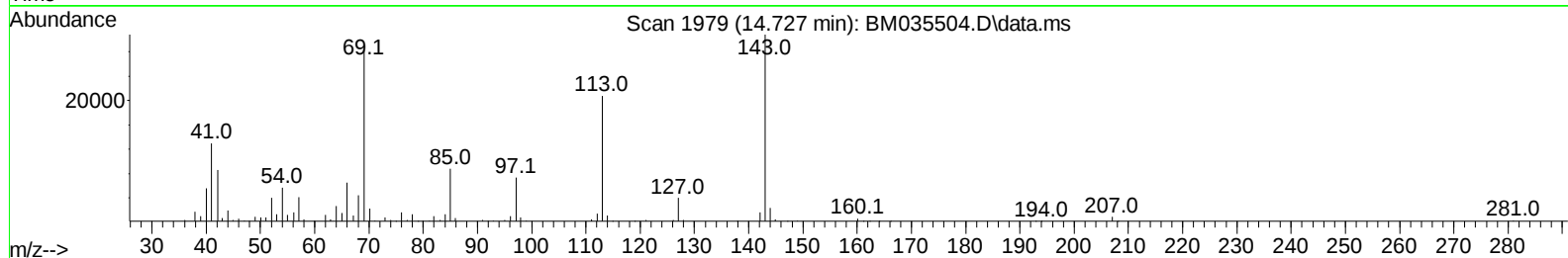
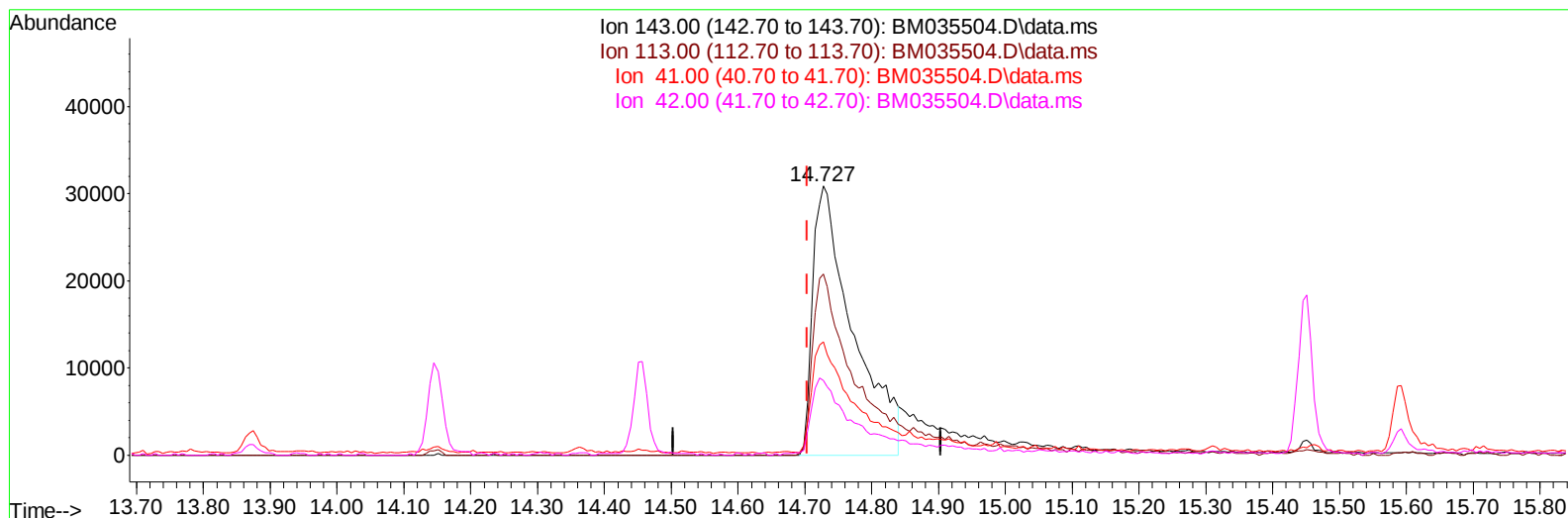
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TIC: BM035504.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.727min (+ 0.024) 21.66 ng/ul m

response 128169

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	61.40	67.31
41.00	41.00	42.11
42.00	29.30	27.74

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.816	152	221594	20.000	ng/ul	0.00
20) Naphthalene-d8	10.610	136	950769	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.457	164	566230	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.204	188	1142699	20.000	ng/ul	0.00
79) Chrysene-d12	21.386	240	1163148	20.000	ng/ul	0.00
88) Perylene-d12	23.727	264	1277098	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.263	96	29992	5.403	ng/uL	0.00
4) Pyridine-d5	3.687	84	251247	18.377	ng/ul	0.00
7) Phenol-d5	6.998	99	498109	29.330	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.145	67	328085	31.066	ng/ul	0.00
11) 2-Chlorophenol-d4	7.351	132	467120	31.756	ng/ul	0.00
15) 4-Methylphenol-d8	8.540	113	391931	27.568	ng/ul	0.00
21) Nitrobenzene-d5	8.975	128	220109	31.101	ng/ul	0.00
24) 2-Nitrophenol-d4	9.698	143	237217	31.084	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.251	165	367624	28.372	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131	549817	26.968	ng/ul	0.00
46) Dimethylphthalate-d6	13.869	166	1279816	33.228	ng/ul	0.00
49) Acenaphthylene-d8	14.145	160	1563161	32.750	ng/ul	0.00
54) 4-Nitrophenol-d4	14.727	143	128169m	21.663	ng/ul	0.02
60) Fluorene-d10	15.451	176	1134148	34.008	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200	154120	22.623	ng/ul	0.00
73) Anthracene-d10	17.304	188	1743056	33.936	ng/ul	0.00
81) Pyrene-d10	19.598	212	2037858	35.637	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.574	264	2166295	35.169	ng/ul	0.00

Target Compounds Qvalue

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

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