

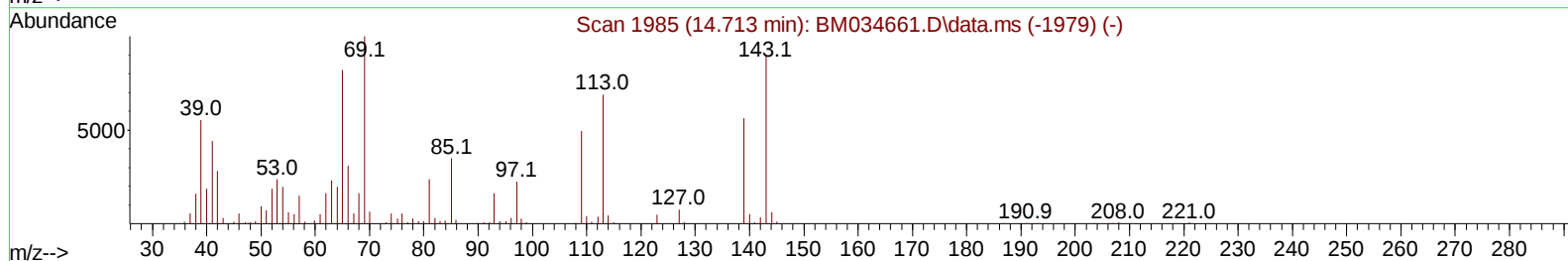
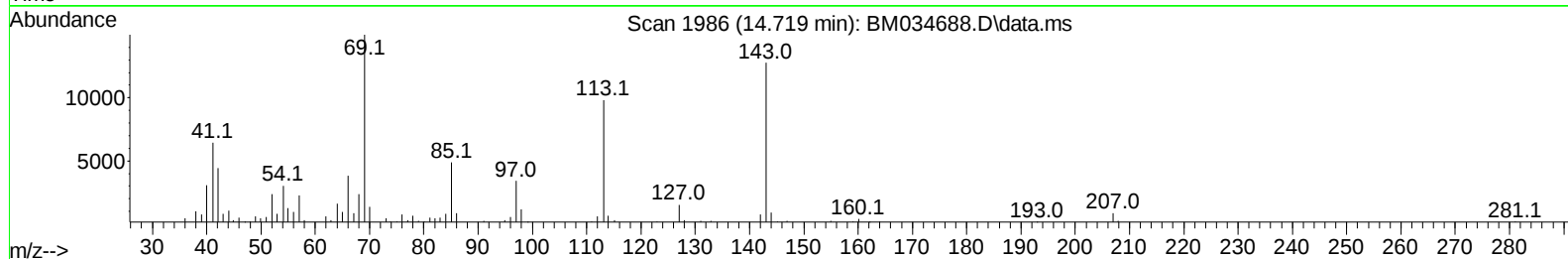
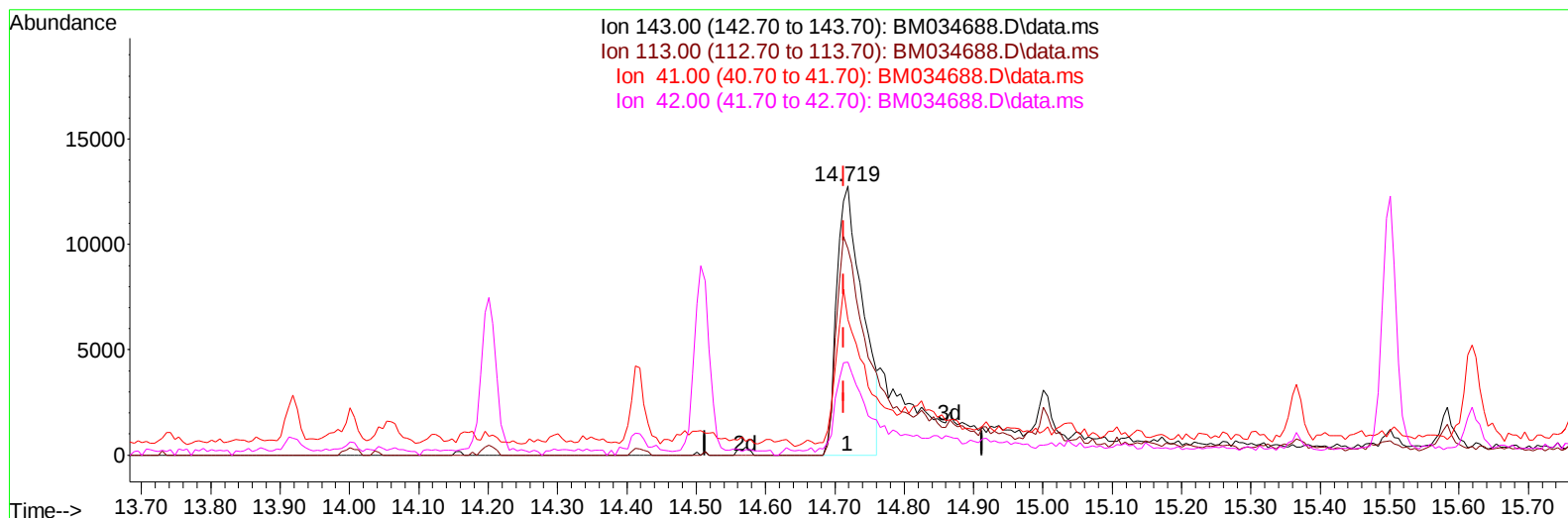
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
 Data File : BM034688.D
 Acq On : 15 Apr 2022 03:39
 Operator : CG/JU
 Sample : N2316-12
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 ETNR2

Manual IntegrationsAPPROVED

Quant Time: Apr 15 04:40:02 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Apr 14 23:42:00 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/15/2022
 Supervised By :Yogesh Patel 04/23/2022



TIC: BM034688.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.719min (+ 0.006) 9.00 ng/u1

response 33093

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	76.83
41.00	31.90	50.38#
42.00	19.10	34.53#

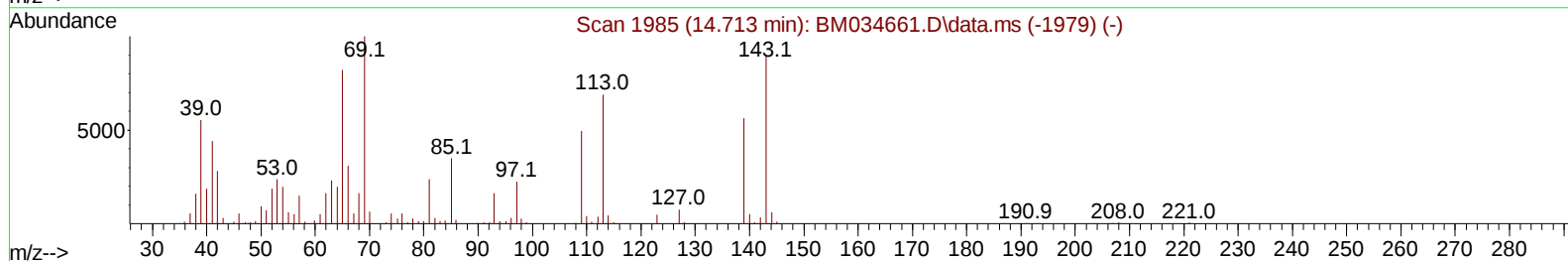
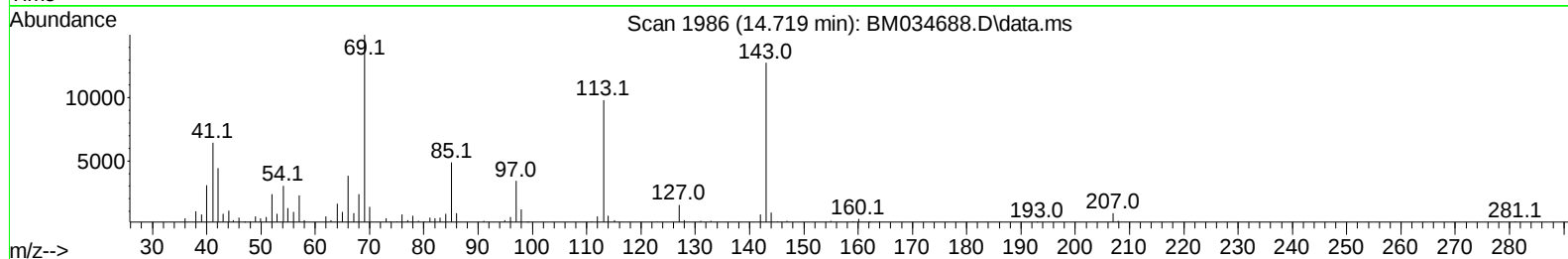
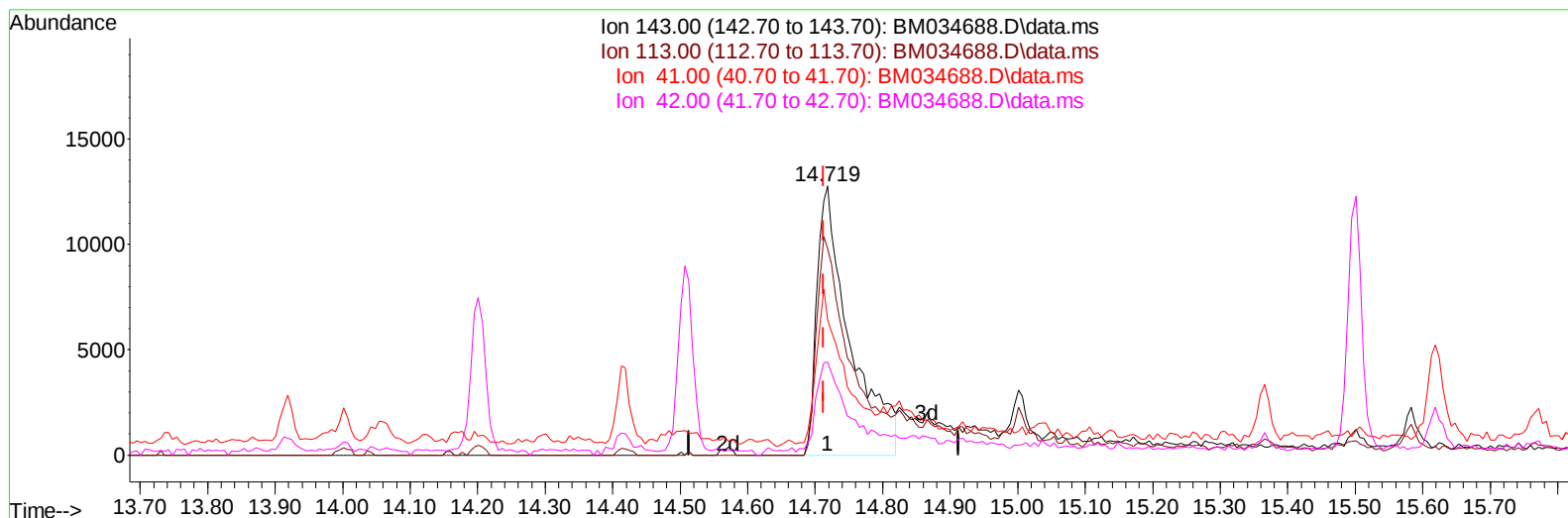
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TIC: BM034688.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.719min (+ 0.006) 11.76 ng/ul m

response 43230

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	76.83
41.00	31.90	50.38#
42.00	19.10	34.53#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.860	152	137316	20.000	ng/ul	0.00
20) Naphthalene-d8	10.672	136	562791	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.507	164	333660	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.254	188	649478	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.430	240	608367	20.000	ng/ul	# 0.00
88) Perylene-d12	23.800	264	657467	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.243	96	16369	4.251	ng/uL	0.00
4) Pyridine-d5	3.666	84	177951	19.149	ng/ul	0.00
7) Phenol-d5	7.025	99	251885	22.039	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.190	67	188138	24.703	ng/ul	0.00
11) 2-Chlorophenol-d4	7.390	132	217916	24.665	ng/ul	0.00
15) 4-Methylphenol-d8	8.572	113	214407	24.016	ng/ul	-0.01
21) Nitrobenzene-d5	9.025	128	102547	24.793	ng/ul	0.00
24) 2-Nitrophenol-d4	9.748	143	105460	24.490	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.289	165	183445	22.623	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.807	131	247832	22.154	ng/ul	0.00
46) Dimethylphthalate-d6	13.919	166	626925	26.003	ng/ul	0.00
49) Acenaphthylene-d8	14.201	160	749312	24.181	ng/ul	0.00
54) 4-Nitrophenol-d4	14.719	143	43230m	11.759	ng/ul	0.00
60) Fluorene-d10	15.501	176	536918	25.770	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.618	200	64133	16.175	ng/ul	0.00
73) Anthracene-d10	17.348	188	791647	25.566	ng/ul	0.00
81) Pyrene-d10	19.642	212	940132	28.313	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.647	264	831797	25.151	ng/ul	-0.01

Target Compounds Qvalue

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

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