

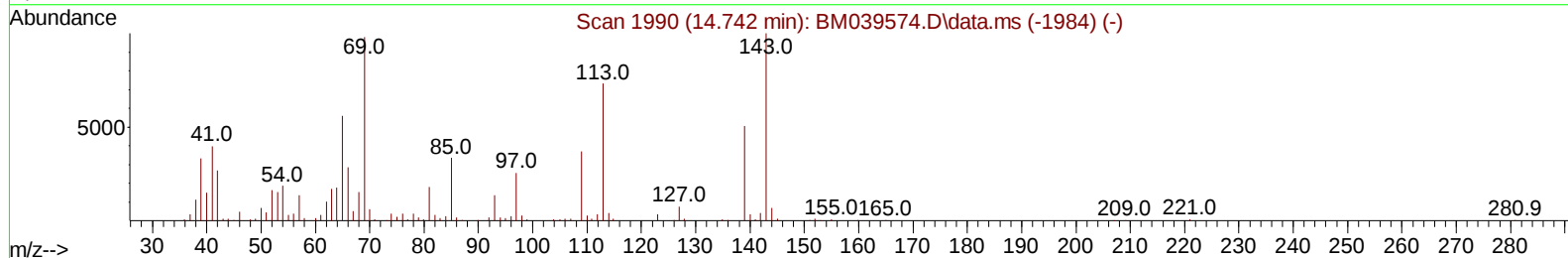
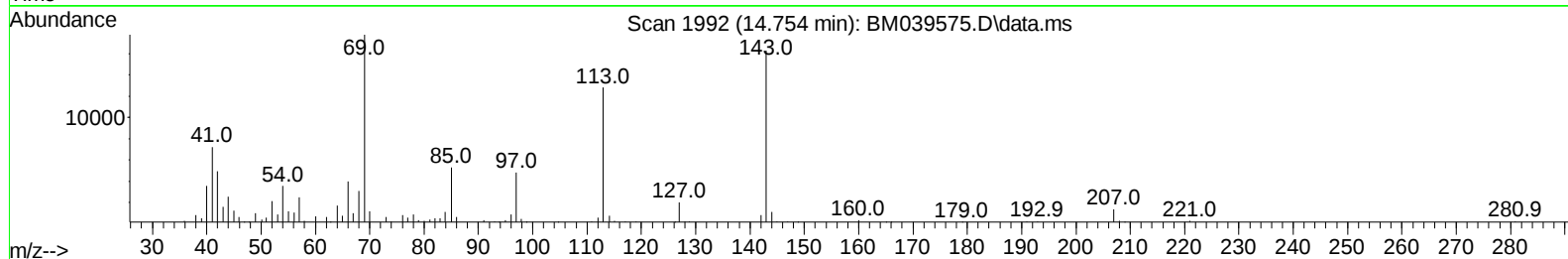
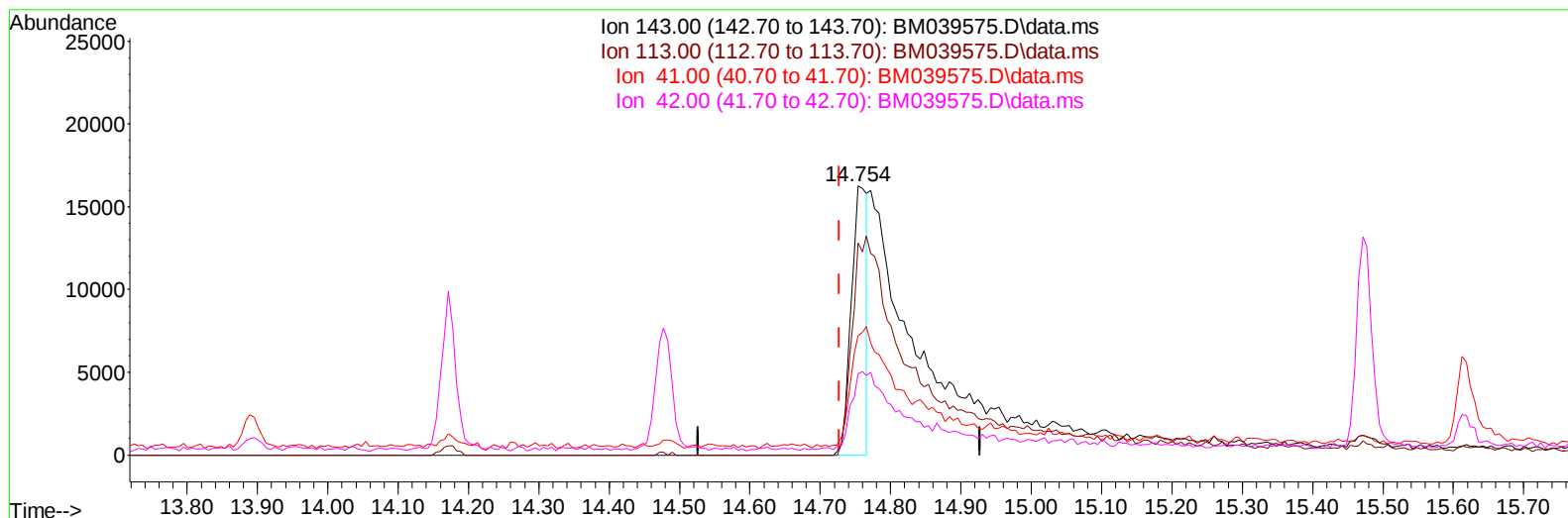
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042223\
Data File : BM039575.D
Acq On : 22 Apr 2023 14:12
Operator : CG/JU
Sample : PB152287BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SBLK287

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/24/2023
Supervised By :Jagrut Upadhyay 04/24/2023

Quant Time: Apr 22 23:43:44 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 19 08:18:41 2023
Response via : Initial Calibration



TIC: BM039575.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.754min (+ 0.026) 5.06 ng/u1

response 25384

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	78.77
41.00	39.90	44.35
42.00	26.60	30.29

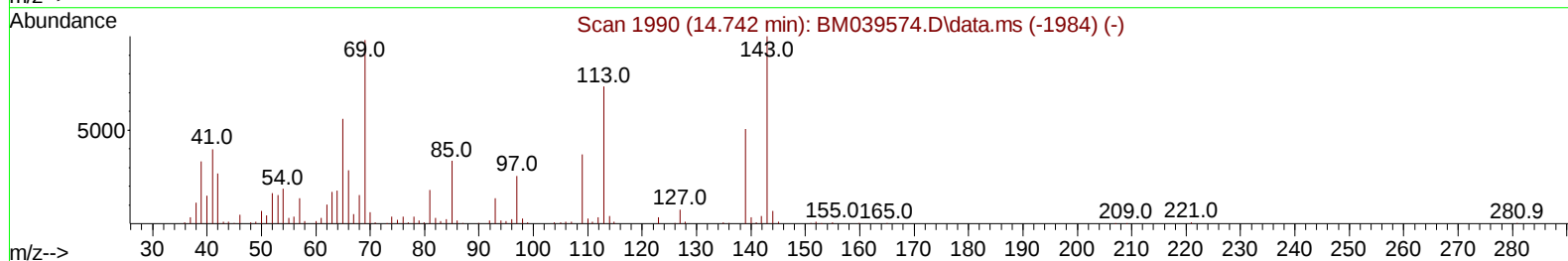
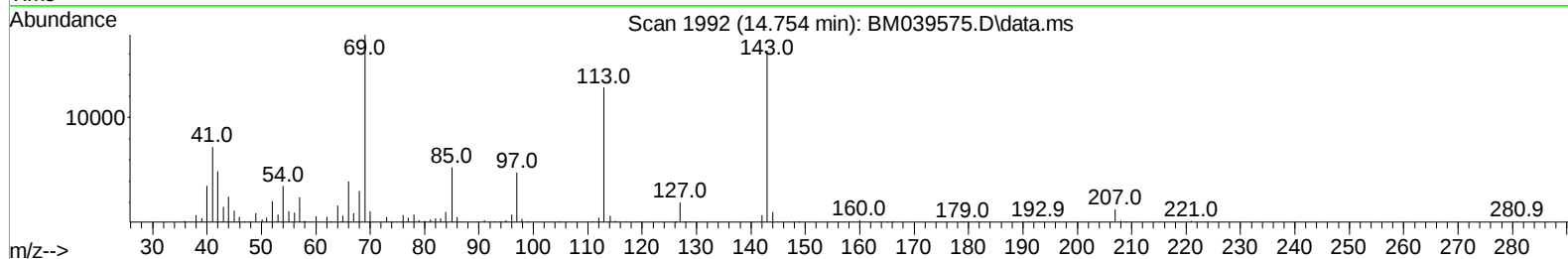
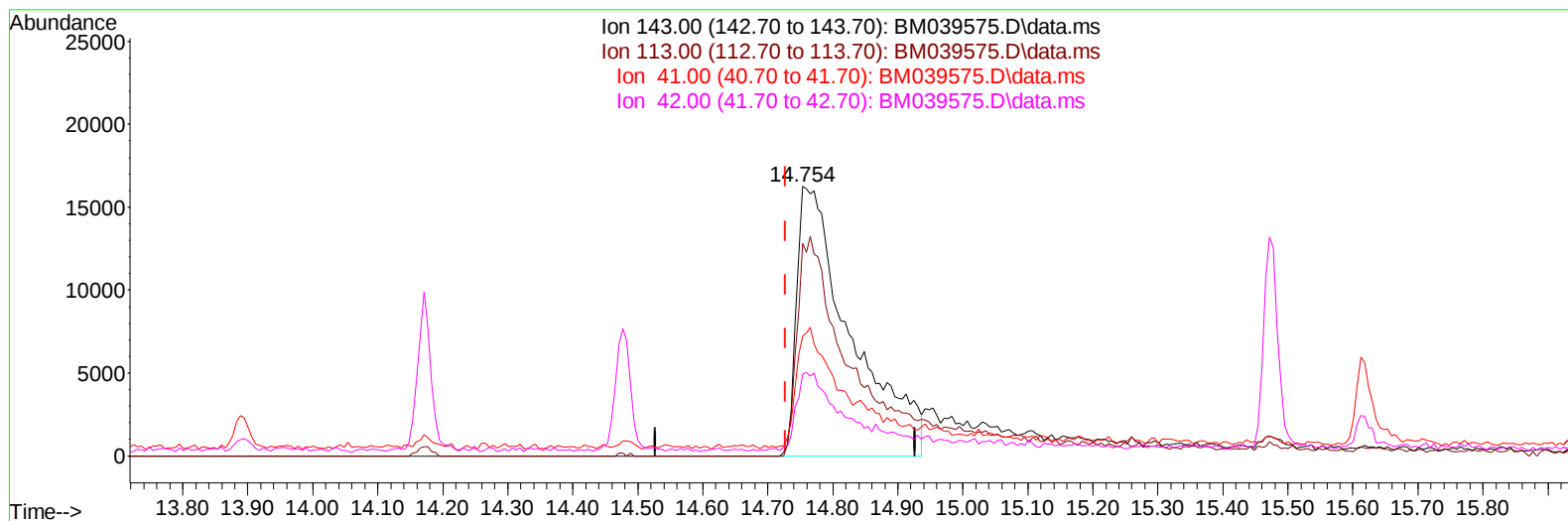
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TIC: BM039575.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.754min (+ 0.026) 18.73 ng/ul m

response 94036

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	78.77
41.00	39.90	44.35
42.00	26.60	30.29

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.819	152	175015	20.000	ng/ul	0.00
20) Naphthalene-d8	10.625	136	707452	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.477	164	411409	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.230	188	836900	20.000	ng/ul	0.00
79) Chrysene-d12	21.424	240	727661	20.000	ng/ul	0.00
88) Perylene-d12	23.794	264	787410	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	28495	6.250	ng/uL	0.00
4) Pyridine-d5	3.637	84	328156	25.614	ng/ul	0.00
7) Phenol-d5	7.007	99	423944	26.466	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.154	67	285548	26.970	ng/ul	0.00
11) 2-Chlorophenol-d4	7.354	132	369355	29.395	ng/ul	0.00
15) 4-Methylphenol-d8	8.554	113	332989	25.807	ng/ul	0.00
21) Nitrobenzene-d5	8.989	128	169587	29.796	ng/ul	0.00
24) 2-Nitrophenol-d4	9.713	143	198338	31.034	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.266	165	290031	26.374	ng/ul	0.00
31) 4-Chloroaniline-d4	10.778	131	417995	25.450	ng/ul	0.00
46) Dimethylphthalate-d6	13.889	166	983417	29.639	ng/ul	-0.01
49) Acenaphthylene-d8	14.171	160	1153266	30.940	ng/ul	0.00
54) 4-Nitrophenol-d4	14.754	143	94036m	18.727	ng/ul	0.03
60) Fluorene-d10	15.471	176	817286	30.065	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.618	200	109259	19.646	ng/ul	0.00
73) Anthracene-d10	17.330	188	1261005	31.331	ng/ul	0.00
81) Pyrene-d10	19.630	212	1416806	29.301	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.641	264	1304849	30.694	ng/ul	0.00

Target Compounds Qvalue

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

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