

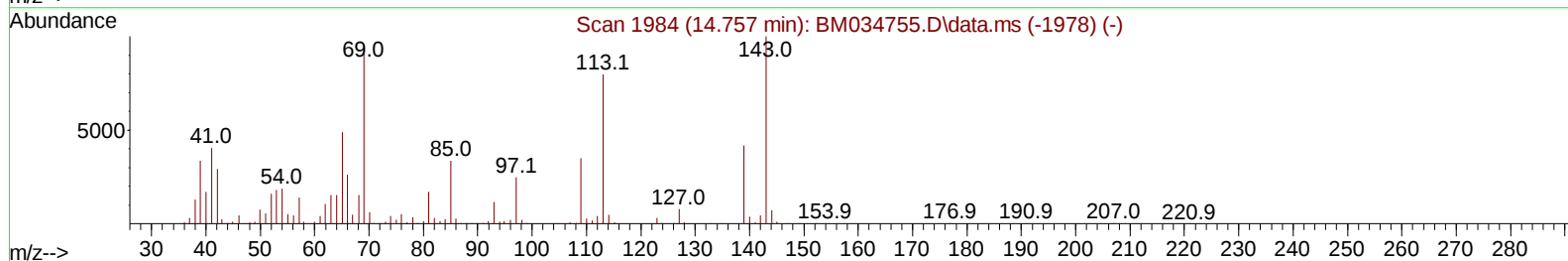
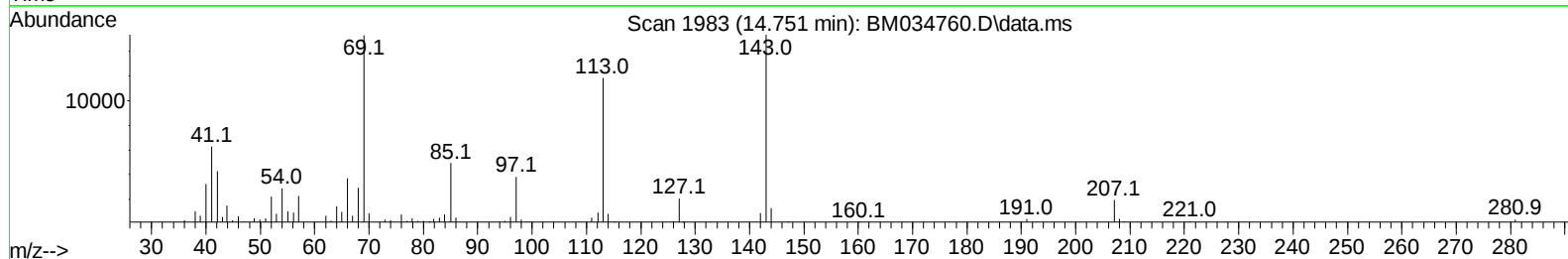
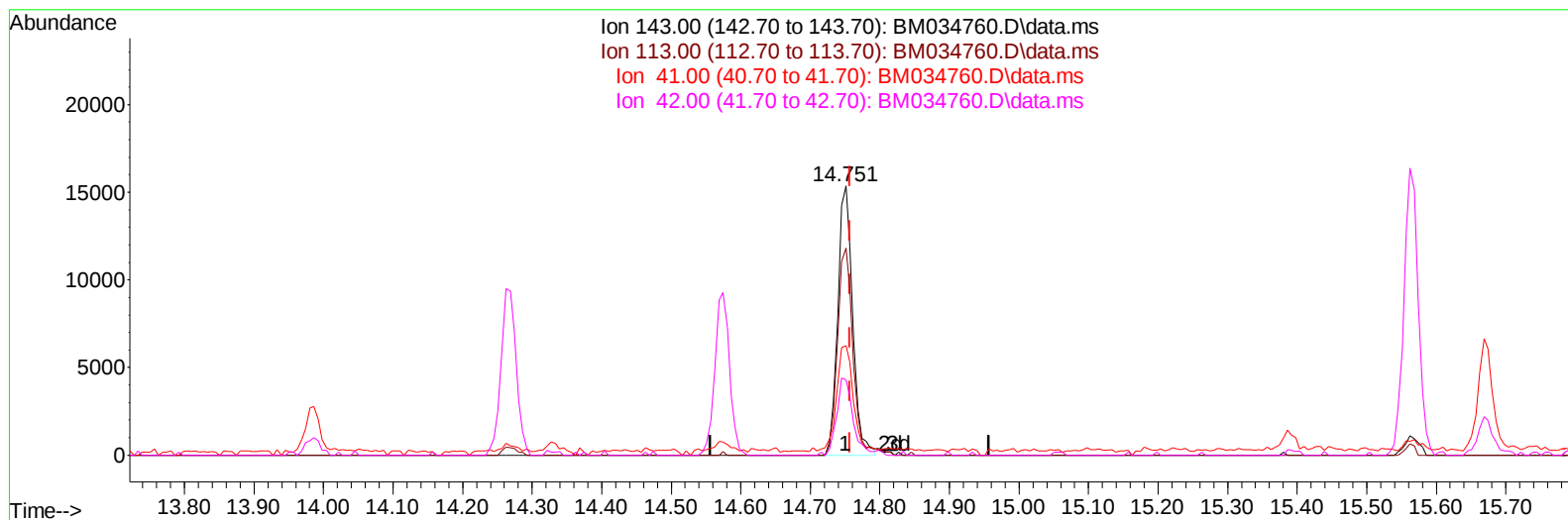
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042122\
Data File : BM034760.D
Acq On : 21 Apr 2022 23:29
Operator : CG/JU
Sample : N2473-19
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
C0J95

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/22/2022
Supervised By :mohammad ahmed 04/26/2022

Quant Time: Apr 22 01:52:40 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM042122.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Apr 21 05:03:15 2022
Response via : Initial Calibration



TIC: BM034760.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.006) 4.72 ng/u1

response 22808

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.00	77.05
41.00	42.20	40.79
42.00	28.30	28.00

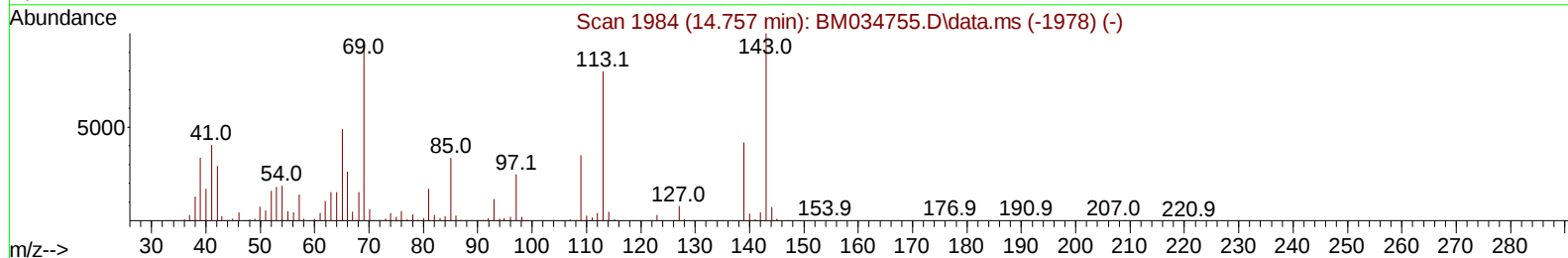
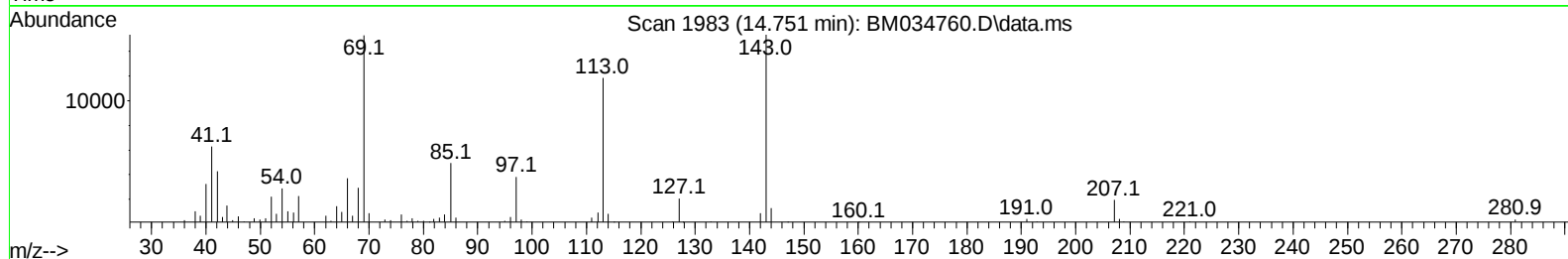
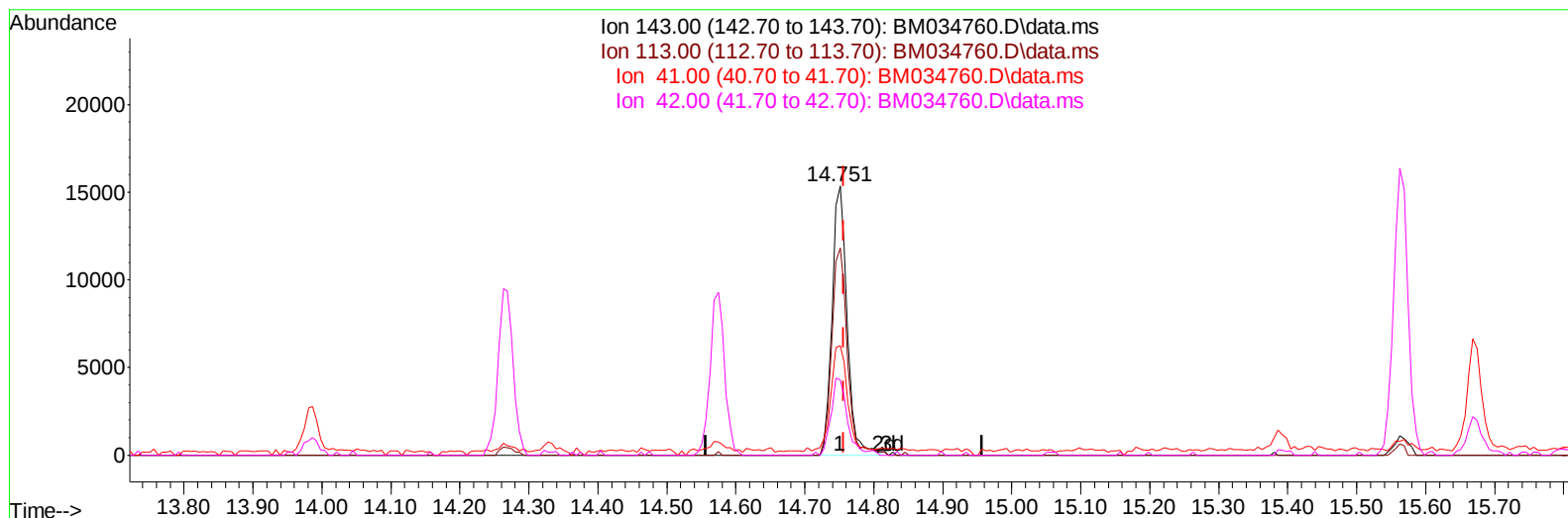
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TIC: BM034760.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.751min (-0.006) 4.77 ng/ul m

response 23062

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.00	77.05
41.00	42.20	40.79
42.00	28.30	28.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.945	152	174618	20.000	ng/ul	0.00
20) Naphthalene-d8	10.745	136	668551	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.574	164	395426	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.315	188	770139	20.000	ng/ul	0.00
79) Chrysene-d12	21.486	240	712366	20.000	ng/ul	-0.01
88) Perylene-d12	23.886	264	744320	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.381	96	21329	5.367	ng/uL	0.00
4) Pyridine-d5	3.799	84	74199	7.522	ng/ul	0.00
7) Phenol-d5	7.098	99	73966	5.608	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.269	67	231500	31.299	ng/ul	0.00
11) 2-Chlorophenol-d4	7.475	132	237555	21.282	ng/ul	0.00
15) 4-Methylphenol-d8	8.645	113	137476	13.028	ng/ul	-0.01
21) Nitrobenzene-d5	9.098	128	146881	30.076	ng/ul	0.00
24) 2-Nitrophenol-d4	9.822	143	130822	25.800	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.363	165	227831	21.330	ng/ul	0.00
31) 4-Chloroaniline-d4	10.875	131	357365	24.744	ng/ul	0.00
46) Dimethylphthalate-d6	13.986	166	895223	31.098	ng/ul	0.00
49) Acenaphthylene-d8	14.269	160	1048235	28.314	ng/ul	0.00
54) 4-Nitrophenol-d4	14.751	143	23062m	4.772	ng/ul	0.00
60) Fluorene-d10	15.563	176	807680	32.425	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.674	200	96024	21.733	ng/ul	0.00
73) Anthracene-d10	17.415	188	1282323	34.194	ng/ul	0.00
81) Pyrene-d10	19.698	212	1396545	36.021	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.733	264	1346652	35.772	ng/ul	0.00

Target Compounds Qvalue

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

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