

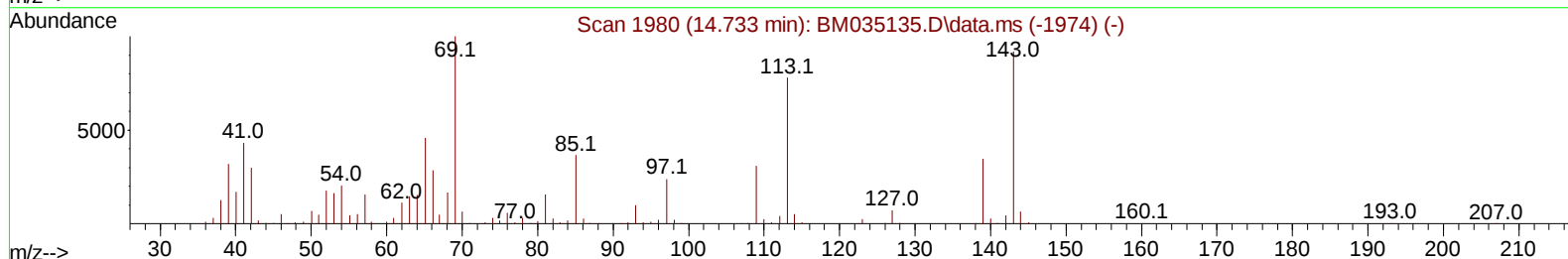
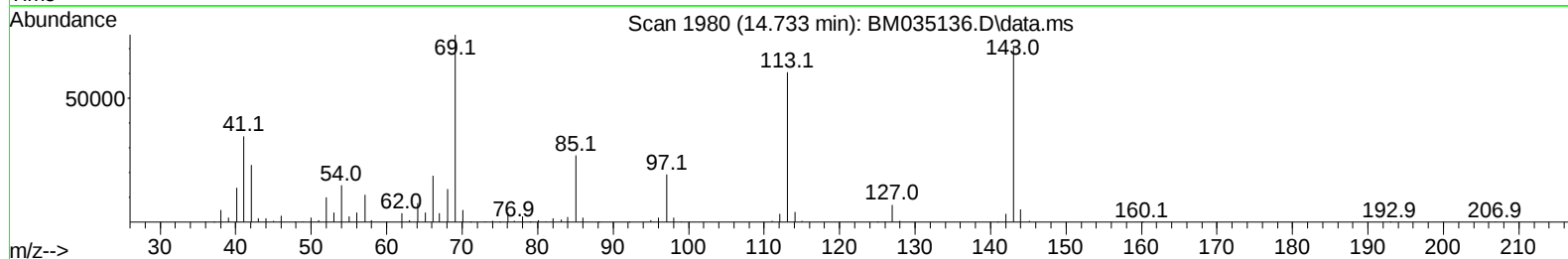
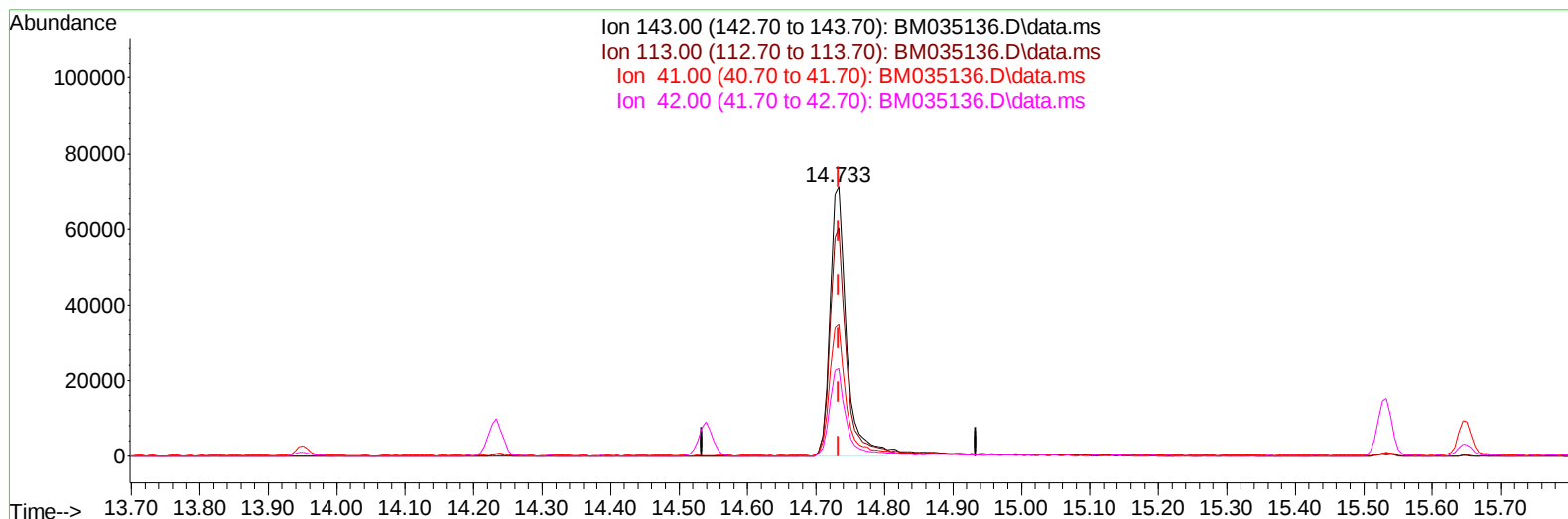
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051822\
 Data File : BM035136.D
 Acq On : 18 May 2022 10:19
 Operator : CG/JU
 Sample : PB144853BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK853

Manual IntegrationsAPPROVED

Quant Time: May 19 00:49:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051722.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue May 17 16:13:52 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/19/2022
 Supervised By :mohammad ahmed 05/19/2022



TIC: BM035136.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (+ 0.000) 27.98 ng/u1

response 119084

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.00	84.63
41.00	42.20	48.70
42.00	28.30	32.42

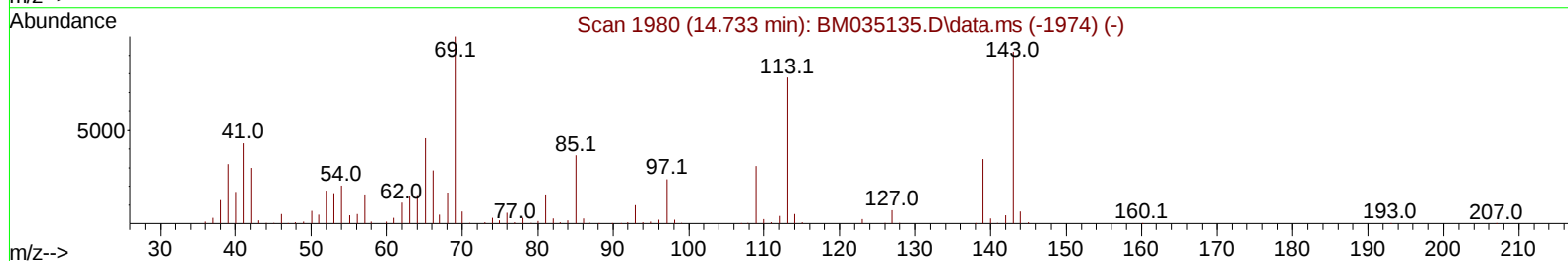
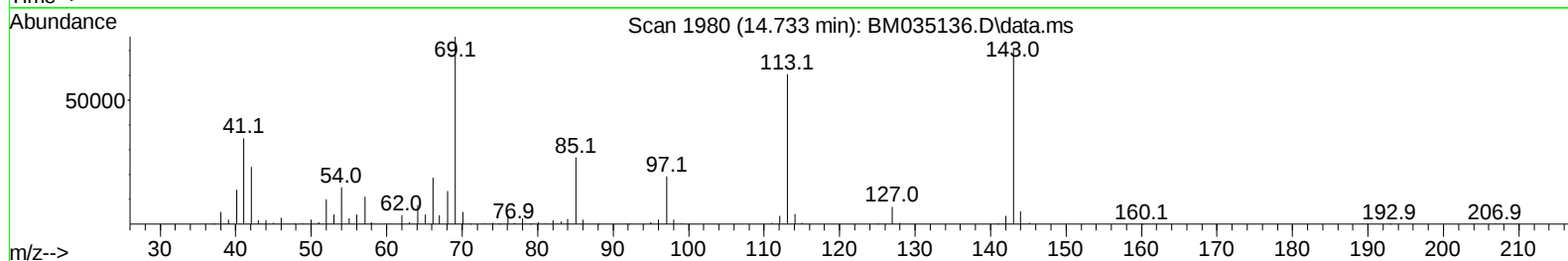
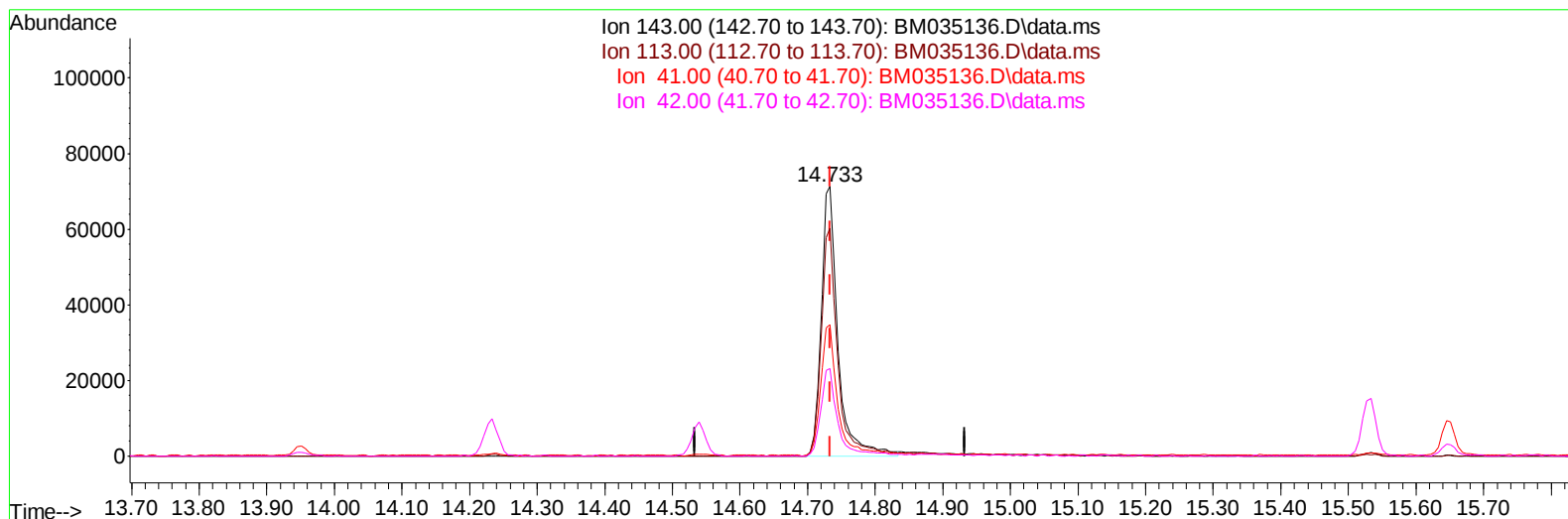
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TIC: BM035136.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.733min (+ 0.000) 28.56 ng/ul m

response 121569

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.00	84.63
41.00	42.20	48.70
42.00	28.30	32.42

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.904	152	144922	20.000	ng/ul	0.00
20) Naphthalene-d8	10.704	136	551382	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.539	164	329506	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.286	188	640667	20.000	ng/ul	0.00
79) Chrysene-d12	21.462	240	583200	20.000	ng/ul	-0.01
88) Perylene-d12	23.850	264	598689	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	23521	6.706	ng/uL	0.00
4) Pyridine-d5	3.781	84	240011	26.527	ng/ul	0.00
7) Phenol-d5	7.069	99	349699	32.742	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.234	67	232292	34.472	ng/ul	0.00
11) 2-Chlorophenol-d4	7.440	132	317208	33.502	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	289100	32.222	ng/ul	0.00
21) Nitrobenzene-d5	9.063	128	144851	33.507	ng/ul	0.00
24) 2-Nitrophenol-d4	9.786	143	158788	32.805	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	265160	30.798	ng/ul	0.00
31) 4-Chloroaniline-d4	10.839	131	264955	22.319	ng/ul	0.00
46) Dimethylphthalate-d6	13.951	166	844186	35.165	ng/ul	0.00
49) Acenaphthylene-d8	14.233	160	980612	34.177	ng/ul	0.00
54) 4-Nitrophenol-d4	14.733	143	121569m	28.563	ng/ul	0.00
60) Fluorene-d10	15.533	176	699483	34.441	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.651	200	120701	29.176	ng/ul	0.00
73) Anthracene-d10	17.380	188	1078016	36.150	ng/ul	0.00
81) Pyrene-d10	19.674	212	1182929	38.083	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.697	264	1137397	37.487	ng/ul	0.00

Target Compounds Qvalue

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

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