

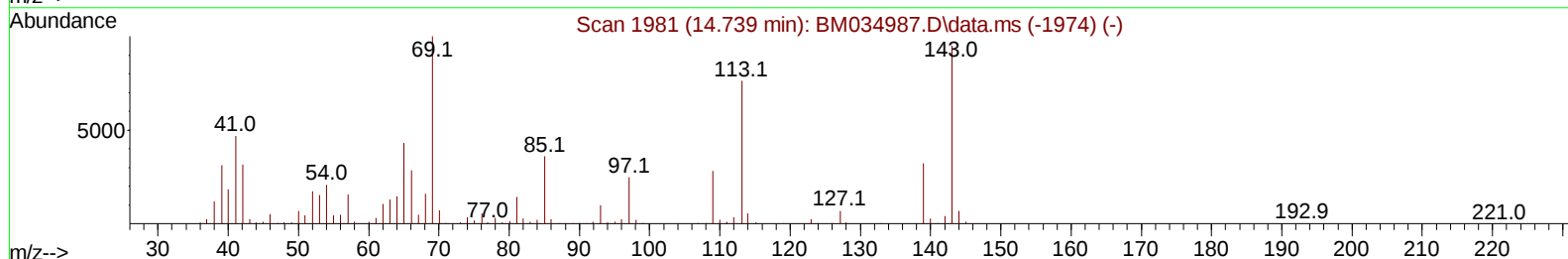
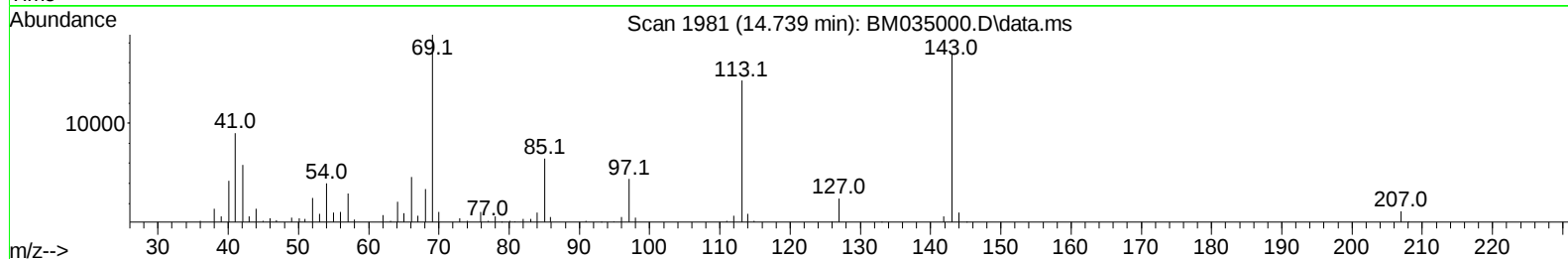
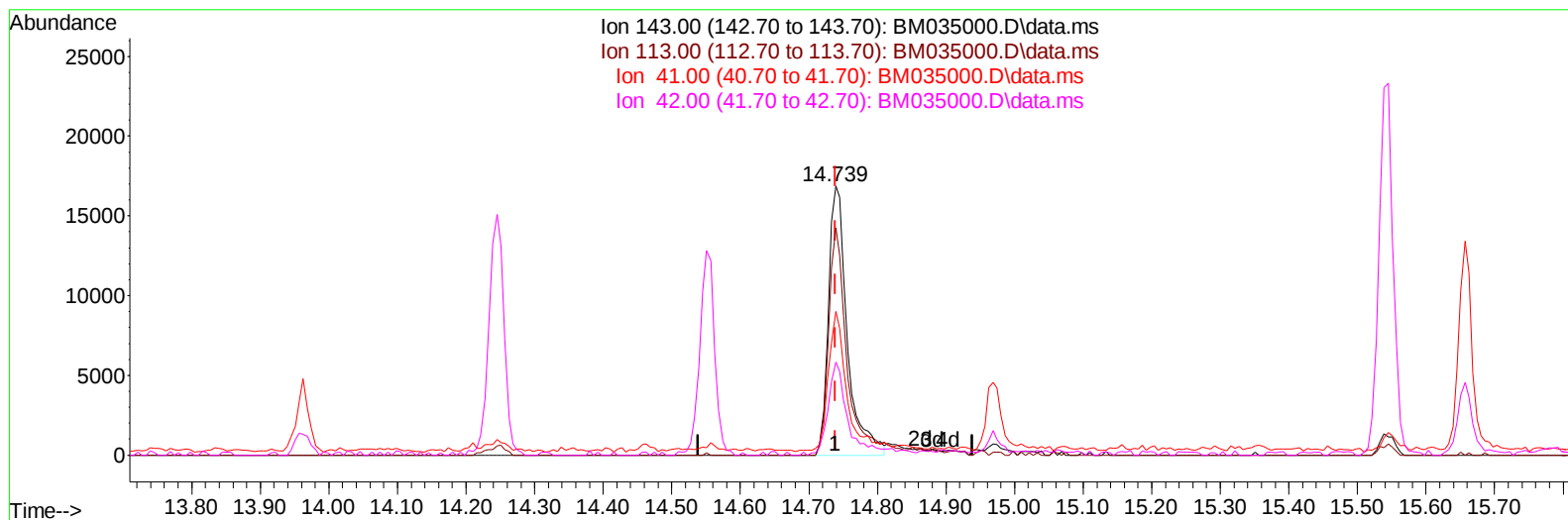
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM051122\
Data File : BM035000.D
Acq On : 12 May 2022 00:10
Operator : CG/JU
Sample : N2707-15
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
EXLA6

Manual IntegrationsAPPROVED

Quant Time: May 12 03:31:21 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM051022.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue May 10 16:40:09 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 05/12/2022
Supervised By :Yogesh Patel 05/17/2022



TIC: BM035000.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (-0.000) 4.86 ng/u1

response 32232

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.00	84.44
41.00	42.20	53.36#
42.00	28.30	34.45#

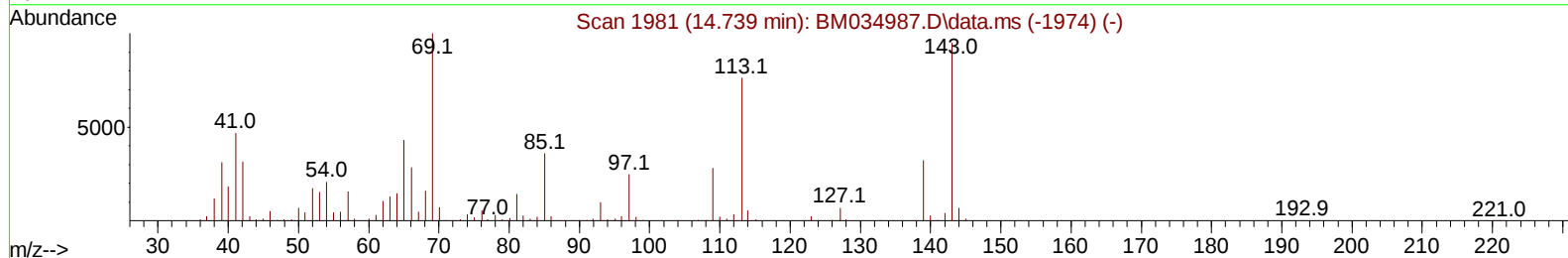
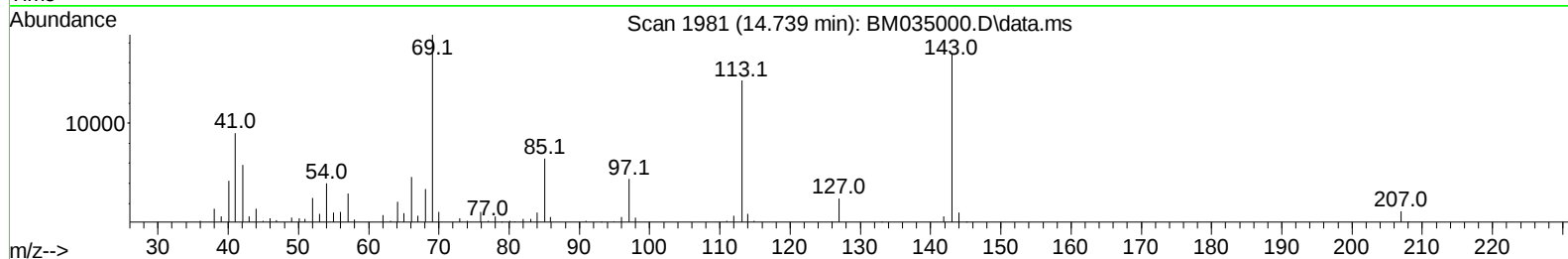
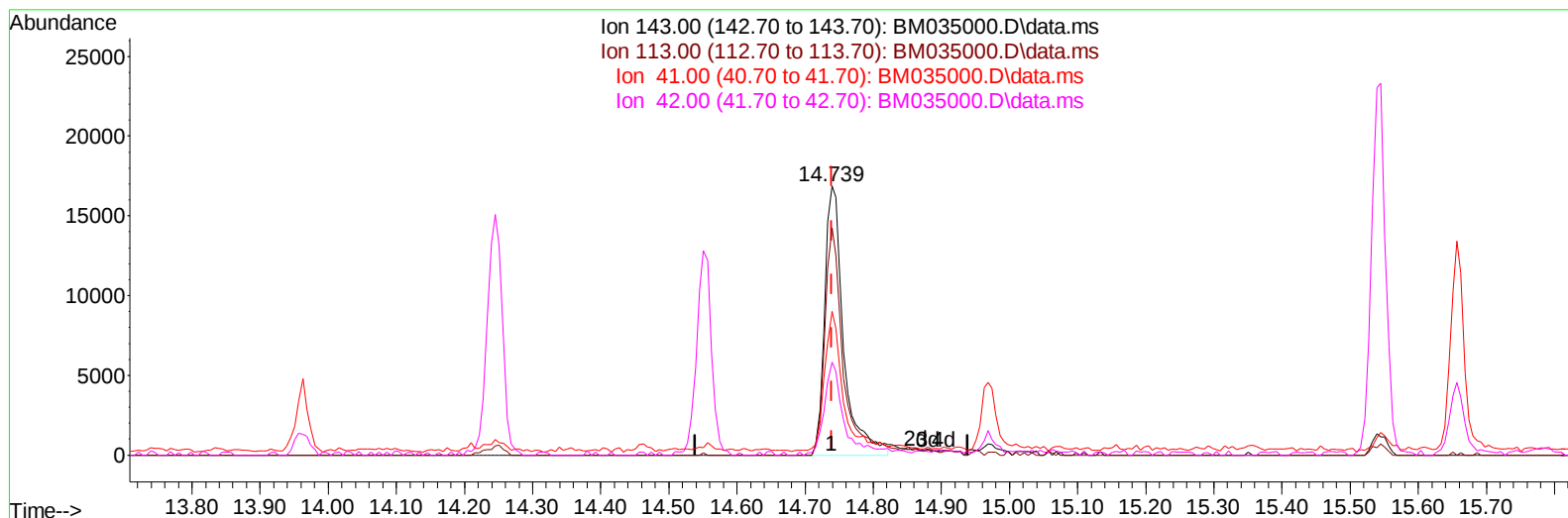
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TIC: BM035000.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.739min (-0.000) 4.93 ng/ul m

response 32718

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	80.00	84.44
41.00	42.20	53.36#
42.00	28.30	34.45#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.922	152	235685	20.000	ng/ul	0.00
20) Naphthalene-d8	10.722	136	943059	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.551	164	514950	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.292	188	916046	20.000	ng/ul	0.00
79) Chrysene-d12	21.468	240	782631	20.000	ng/ul	0.00
88) Perylene-d12	23.850	264	878814	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.369	96	22706	4.031	ng/uL	0.00
4) Pyridine-d5	3.781	84	95962	6.180	ng/ul	0.00
7) Phenol-d5	7.081	99	114316	5.607	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.251	67	368183	28.770	ng/ul	0.00
11) 2-Chlorophenol-d4	7.451	132	384083	23.972	ng/ul	0.00
15) 4-Methylphenol-d8	8.628	113	229908	13.830	ng/ul	0.00
21) Nitrobenzene-d5	9.075	128	240602	33.288	ng/ul	0.00
24) 2-Nitrophenol-d4	9.804	143	253149	32.472	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.339	165	395640	27.117	ng/ul	0.00
31) 4-Chloroaniline-d4	10.851	131	499115	24.026	ng/ul	0.00
46) Dimethylphthalate-d6	13.963	166	1309309	33.714	ng/ul	0.00
49) Acenaphthylene-d8	14.245	160	1566152	31.782	ng/ul	0.00
54) 4-Nitrophenol-d4	14.739	143	32718m	4.931	ng/ul	0.00
60) Fluorene-d10	15.545	176	1091411	33.311	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.657	200	159969	27.712	ng/ul	0.00
73) Anthracene-d10	17.392	188	1586188	35.590	ng/ul	0.00
81) Pyrene-d10	19.680	212	1677641	34.329	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.703	264	1743632	37.963	ng/ul	0.00

Target Compounds Qvalue

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

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