

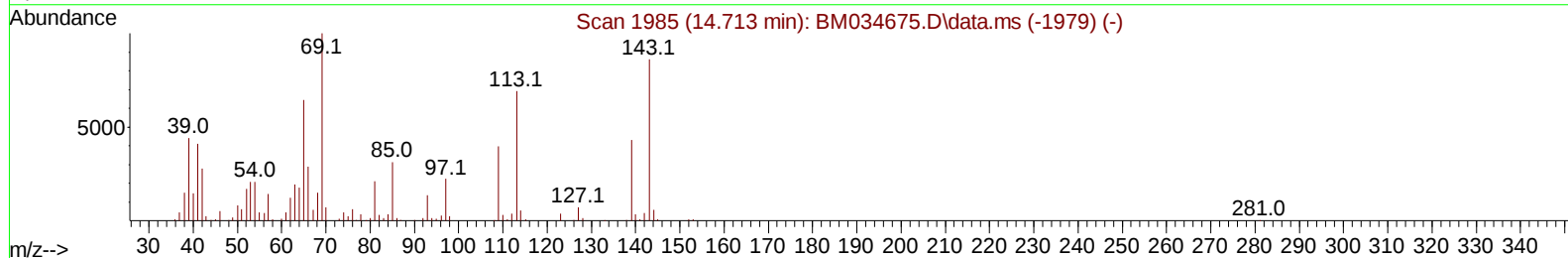
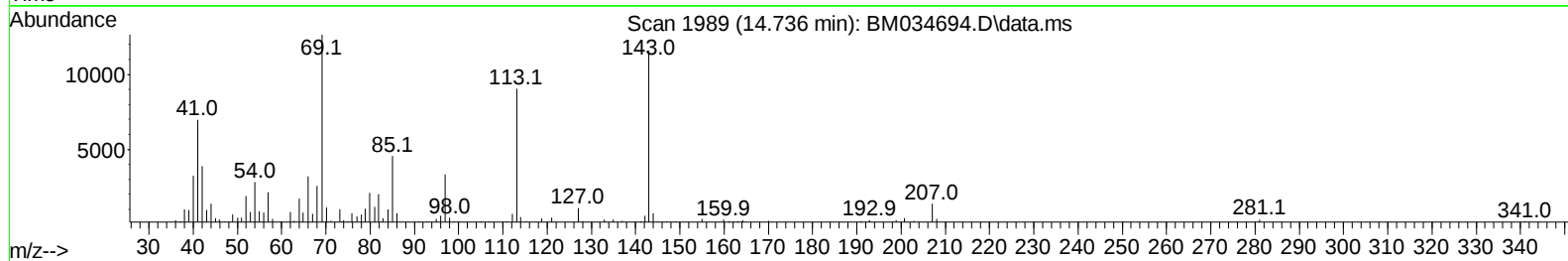
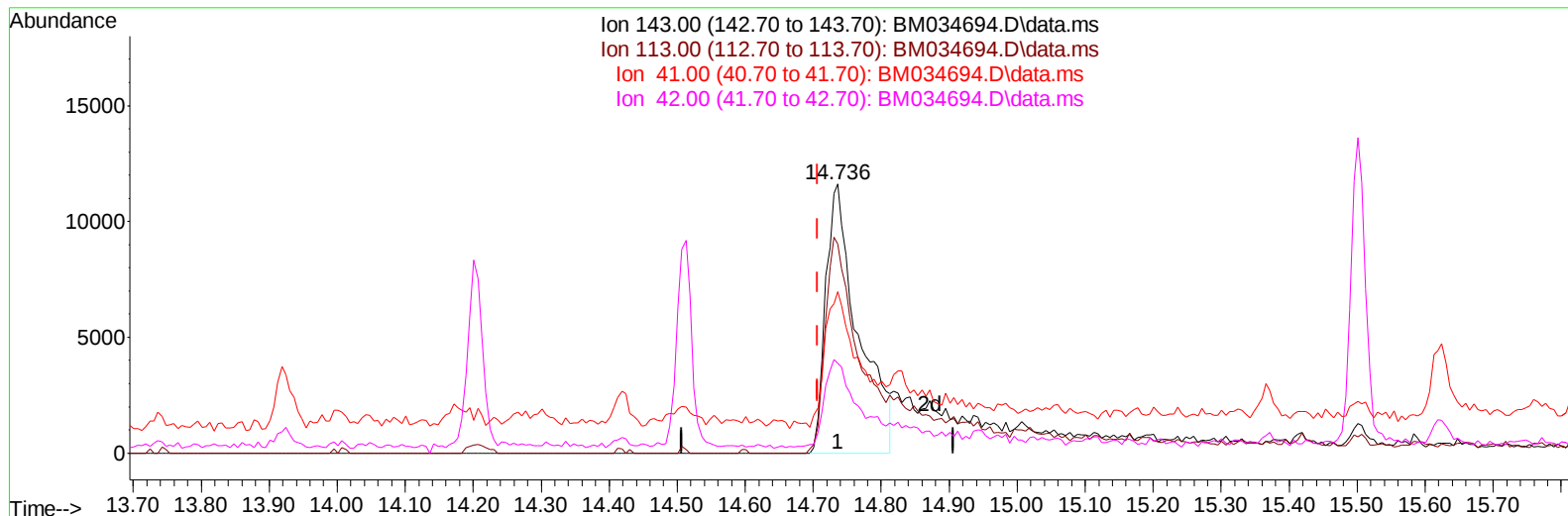
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM041422\
Data File : BM034694.D
Acq On : 15 Apr 2022 07:51
Operator : CG/JU
Sample : N2316-10
Misc :
ALS Vial : 37 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ETNQ9

Manual IntegrationsAPPROVED

Quant Time: Apr 18 06:31:31 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Apr 18 05:39:41 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 04/22/2022
Supervised By :Yogesh Patel 04/23/2022



TIC: BM034694.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.736min (+ 0.029) 10.08 ng/u1

response 38501

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	77.80
41.00	31.90	60.04#
42.00	19.10	33.35#

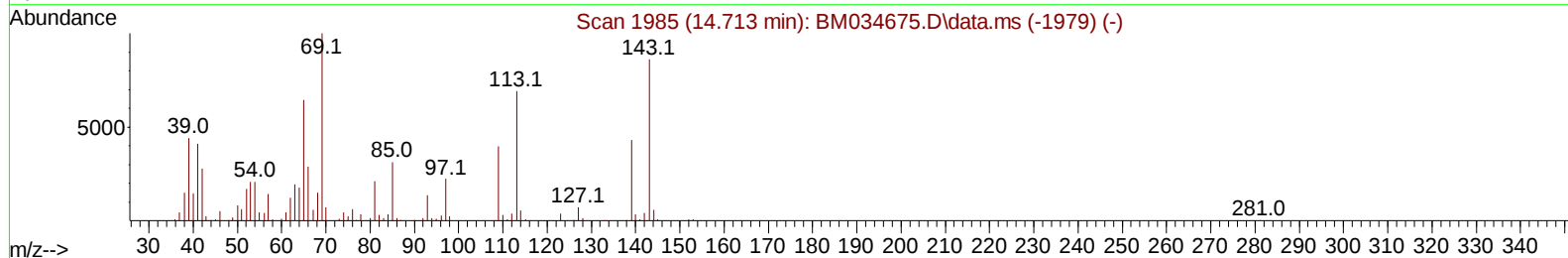
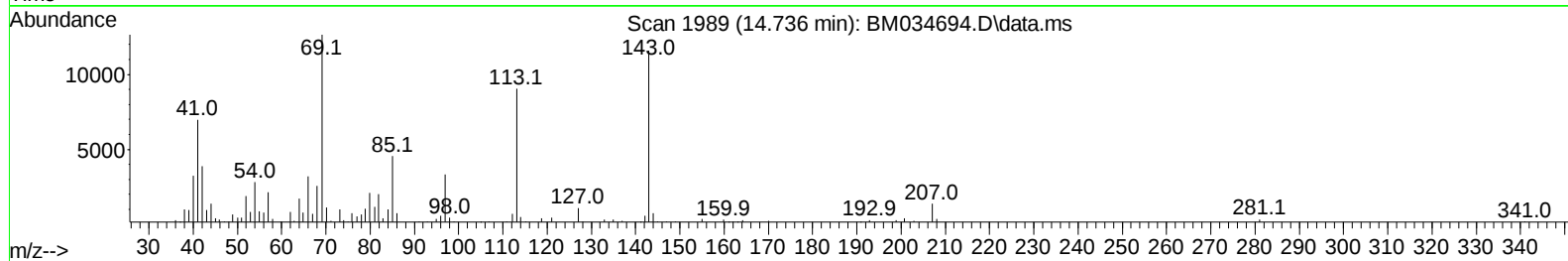
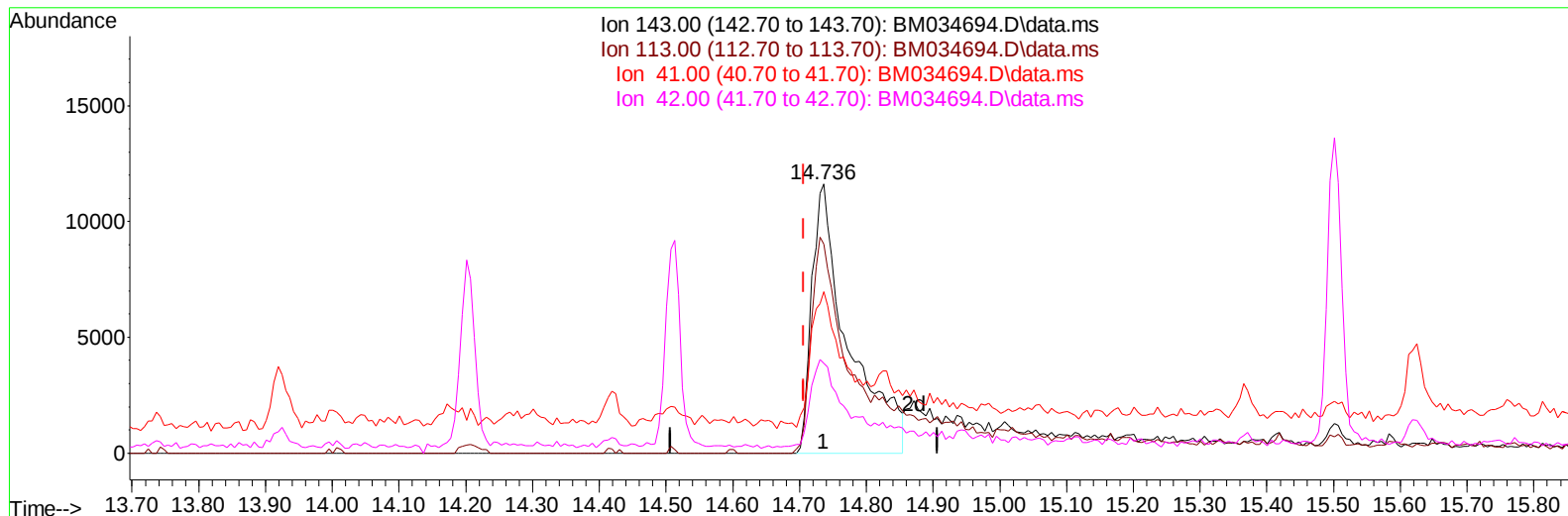
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(54) 4-Nitrophenol-d4 (S)

14.736min (+ 0.029) 11.54 ng/ul m

response 44067

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	77.80
41.00	31.90	60.04#
42.00	19.10	33.35#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.860	152	144286	20.000	ng/ul	0.00
20) Naphthalene-d8	10.672	136	590607	20.000	ng/ul	# 0.01
38) Acenaphthene-d10	14.513	164	346721	20.000	ng/ul	0.01
64) Phenanthrene-d10	17.254	188	694497	20.000	ng/ul	# 0.01
79) Chrysene-d12	21.430	240	562524	20.000	ng/ul	# 0.00
88) Perylene-d12	23.806	264	436093	20.000	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.243	96	15455	3.820	ng/uL	0.01
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.031	99	273615	22.784	ng/ul	0.01
9) Bis-(2-Chloroethyl)eth...	7.190	67	202321	25.282	ng/ul	0.01
11) 2-Chlorophenol-d4	7.396	132	233882	25.193	ng/ul	0.01
15) 4-Methylphenol-d8	8.578	113	227112	24.210	ng/ul	0.01
21) Nitrobenzene-d5	9.025	128	109257	25.171	ng/ul	0.01
24) 2-Nitrophenol-d4	9.754	143	102790	22.745	ng/ul	0.01
28) 2,4-Dichlorophenol-d3	10.295	165	197550	23.215	ng/ul	0.01
31) 4-Chloroaniline-d4	10.819	131	161776	13.780	ng/ul	0.02
46) Dimethylphthalate-d6	13.925	166	720098	28.743	ng/ul	0.01
49) Acenaphthylene-d8	14.201	160	817112	25.376	ng/ul	0.00
54) 4-Nitrophenol-d4	14.736	143	44067m	11.536	ng/ul	0.03
60) Fluorene-d10	15.501	176	617148	28.505	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.624	200	44089	10.399	ng/ul	0.01
73) Anthracene-d10	17.354	188	983726	29.710	ng/ul	0.01
81) Pyrene-d10	19.642	212	1197378	39.000	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.647	264	626617	28.565	ng/ul	0.00

Target Compounds Qvalue

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

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