

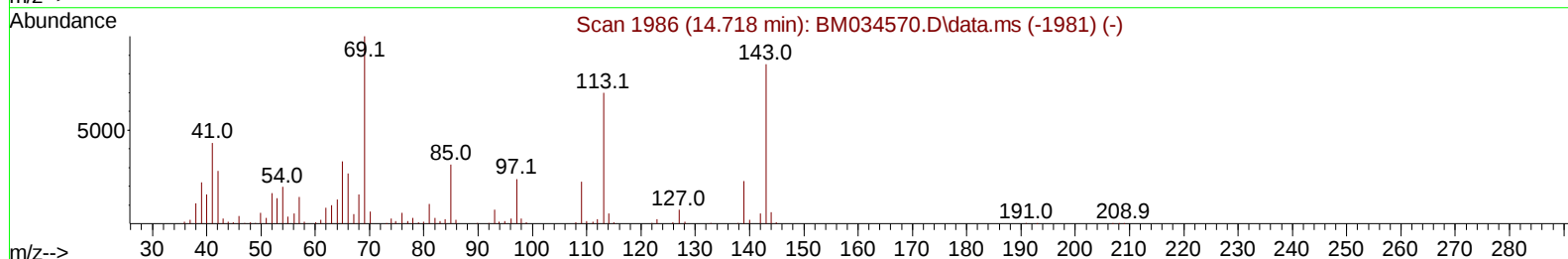
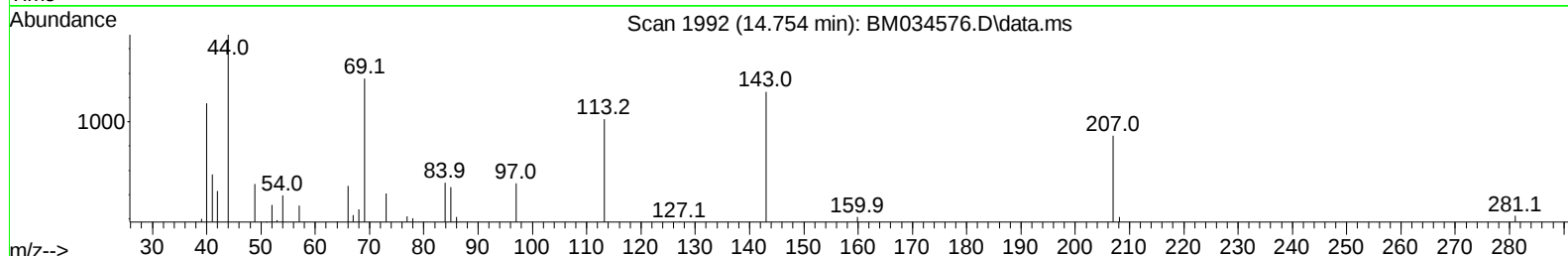
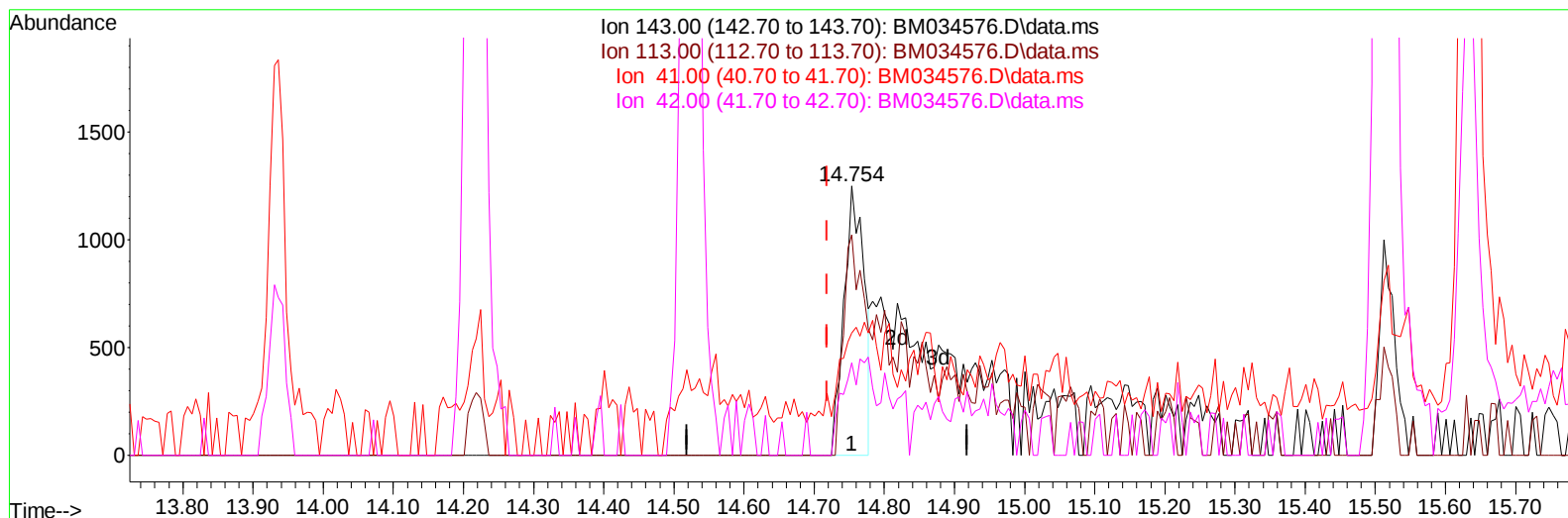
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040922\
Data File : BM034576.D
Acq On : 09 Apr 2022 00:37
Operator : CG/JU
Sample : N2319-09
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
YB2H7

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 04/11/2022
Supervised By :Yogesh Patel 04/12/2022

Quant Time: Apr 09 04:46:39 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM040922.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 08 14:30:08 2022
Response via : Initial Calibration



TIC: BM034576.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.754min (+ 0.035) 0.85 ng/u1

response 2507

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	81.83
41.00	31.90	45.32#
42.00	19.10	34.35#

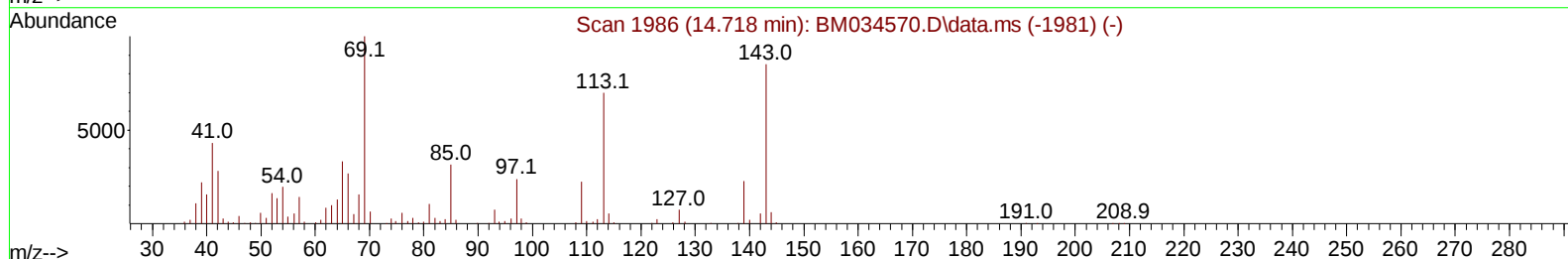
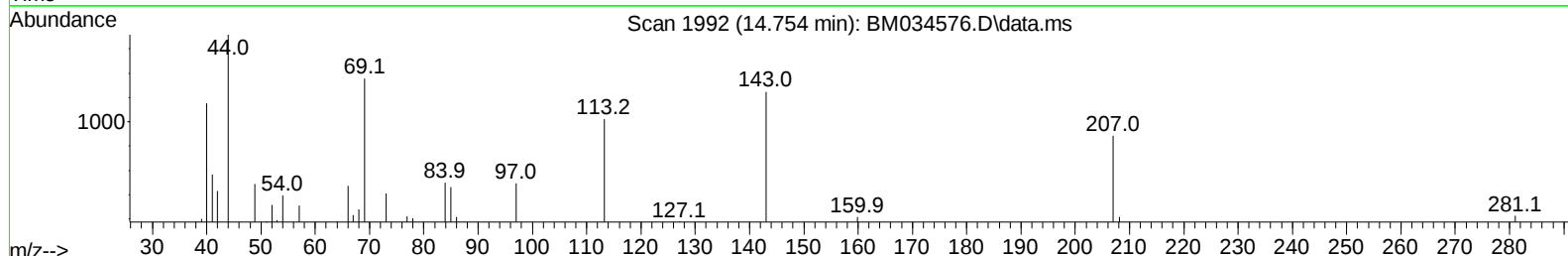
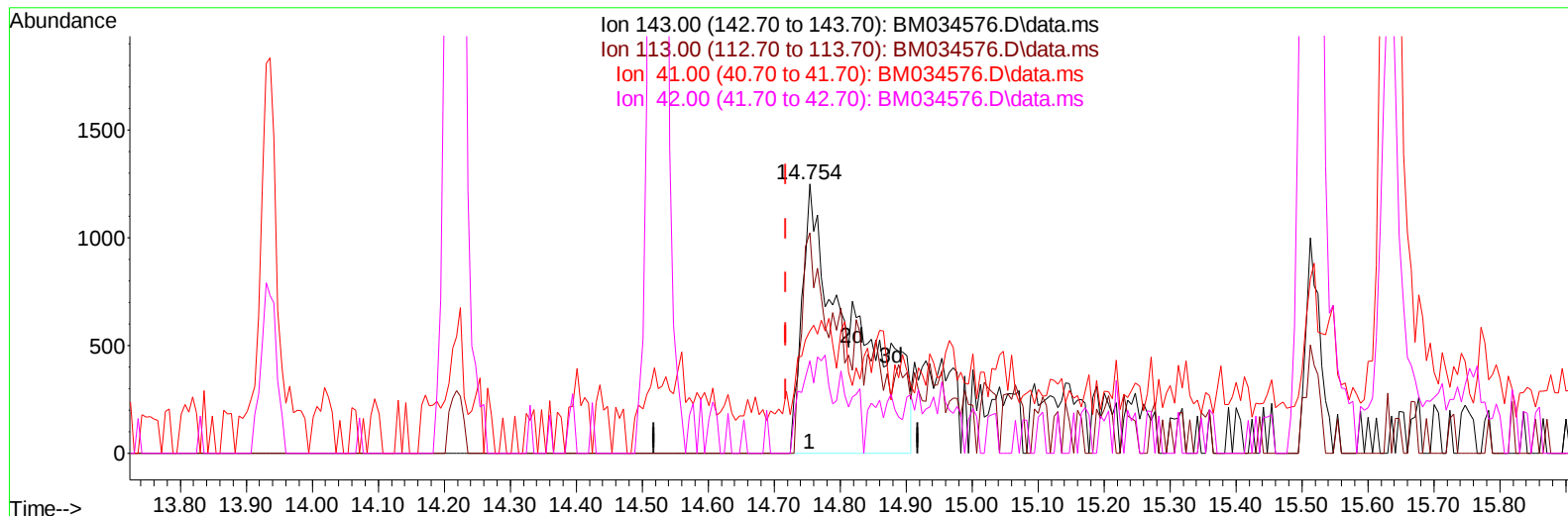
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TIC: BM034576.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.754min (+ 0.035) 2.27 ng/ul m

response 6682

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	81.83
41.00	31.90	45.32#
42.00	19.10	34.35#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.878	152	117371	20.000	ng/ul	0.00
20) Naphthalene-d8	10.683	136	471348	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.524	164	267016	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.265	188	516883	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.442	240	477541	20.000	ng/ul	# 0.00
88) Perylene-d12	23.818	264	537315	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.254	96	13760	4.181	ng/uL	0.00
4) Pyridine-d5	3.690	84	50192	6.319	ng/ul	0.01
7) Phenol-d5	7.042	99	53281	5.454	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.201	67	216501	33.258	ng/ul	0.00
11) 2-Chlorophenol-d4	7.407	132	191924	25.414	ng/ul	0.00
15) 4-Methylphenol-d8	8.589	113	109162	14.305	ng/ul	0.00
21) Nitrobenzene-d5	9.042	128	116190	33.541	ng/ul	0.00
24) 2-Nitrophenol-d4	9.766	143	114781	31.825	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.307	165	181693	26.754	ng/ul	0.00
31) 4-Chloroaniline-d4	10.825	131	234981	25.080	ng/ul	0.00
46) Dimethylphthalate-d6	13.936	166	686343	35.573	ng/ul	0.00
49) Acenaphthylene-d8	14.218	160	831894	33.546	ng/ul	0.00
54) 4-Nitrophenol-d4	14.754	143	6682m	2.271	ng/ul	0.04
60) Fluorene-d10	15.512	176	617417	37.031	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.636	200	93278	29.561	ng/ul	0.00
73) Anthracene-d10	17.365	188	976411	39.622	ng/ul	0.00
81) Pyrene-d10	19.653	212	1173566	45.026	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.665	264	1042926	38.587	ng/ul	0.00

Target Compounds Qvalue

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

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