

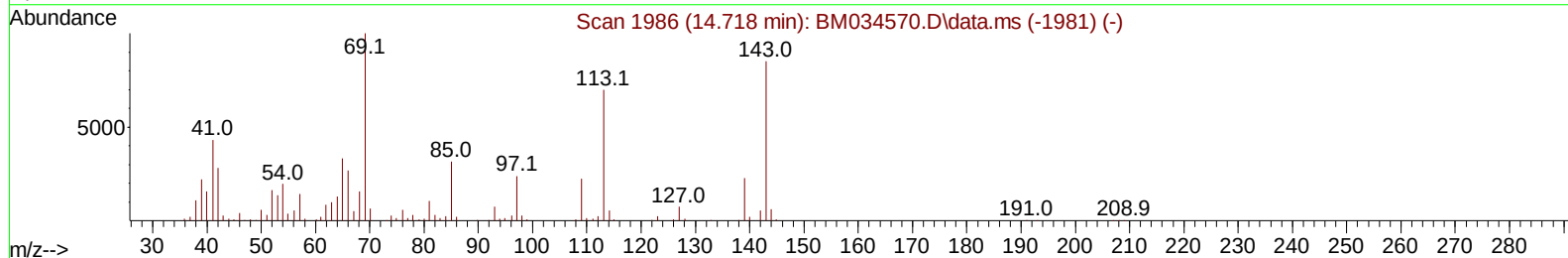
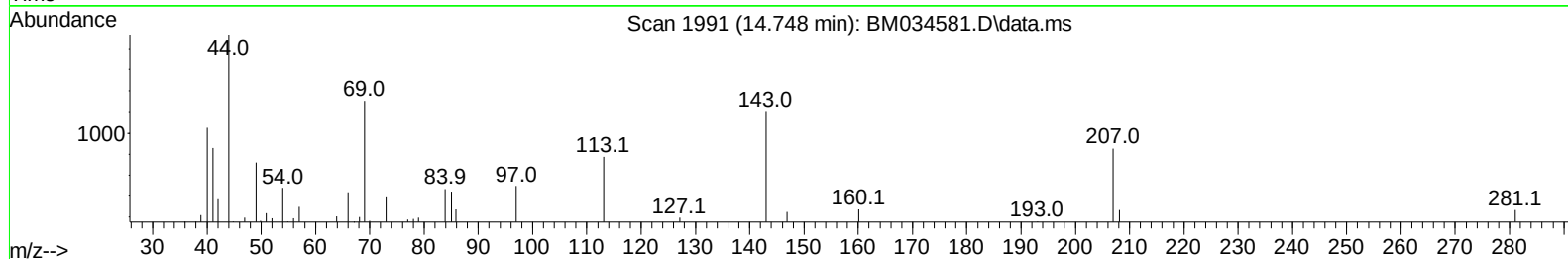
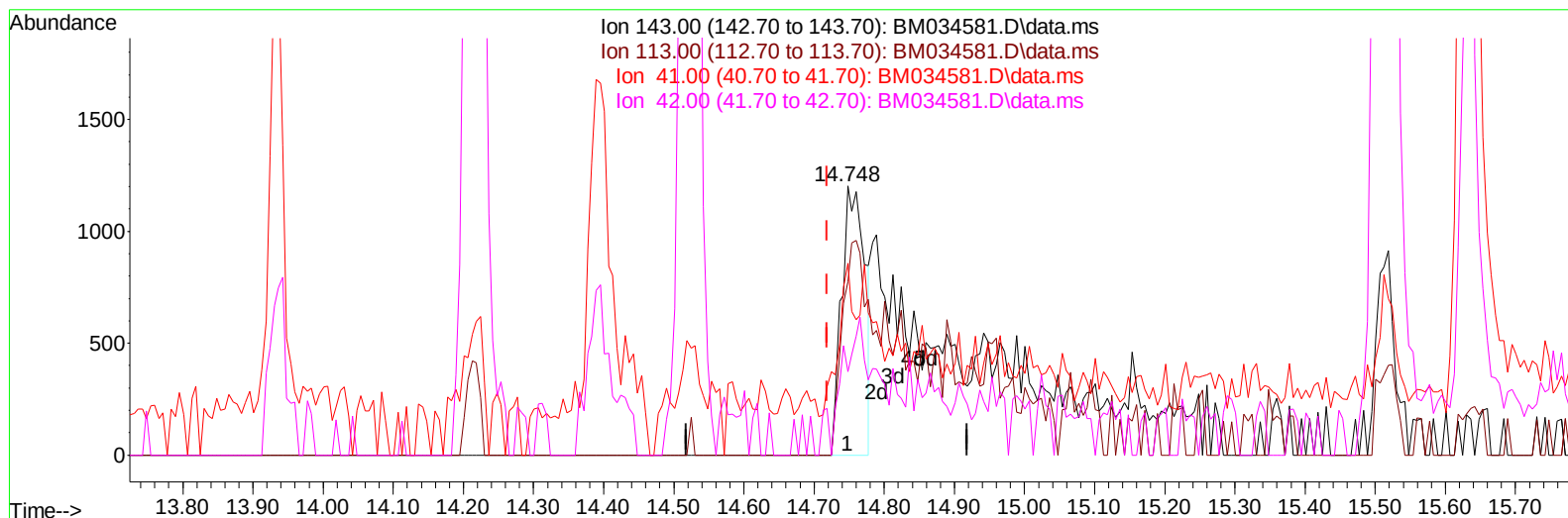
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM040922\
Data File : BM034581.D
Acq On : 09 Apr 2022 03:38
Operator : CG/JU
Sample : N2319-01
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
YB2G7

Manual IntegrationsAPPROVED

Quant Time: Apr 09 04:47:51 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

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



TIC: BM034581.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.748min (+ 0.030) 0.80 ng/ul

response 2767

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	64.48
41.00	31.90	71.30#
42.00	19.10	31.03#

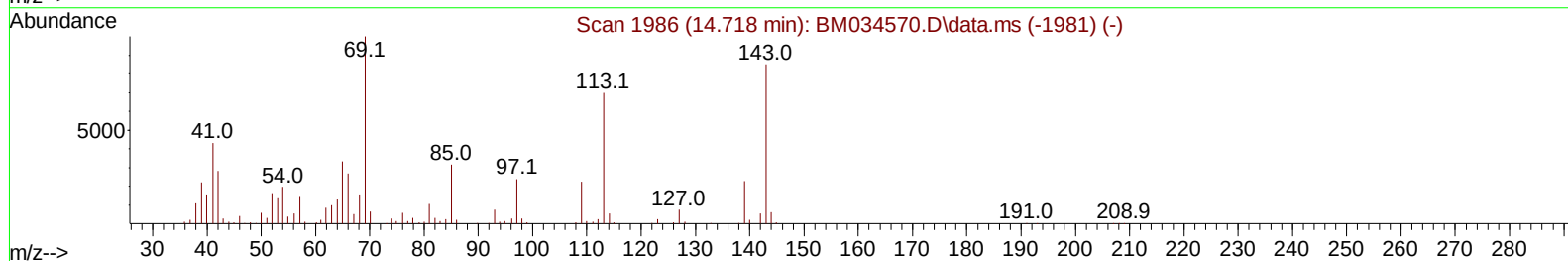
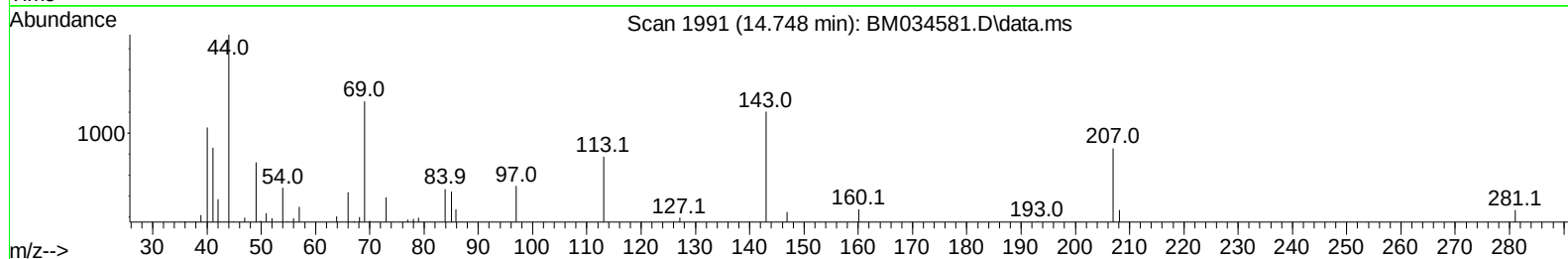
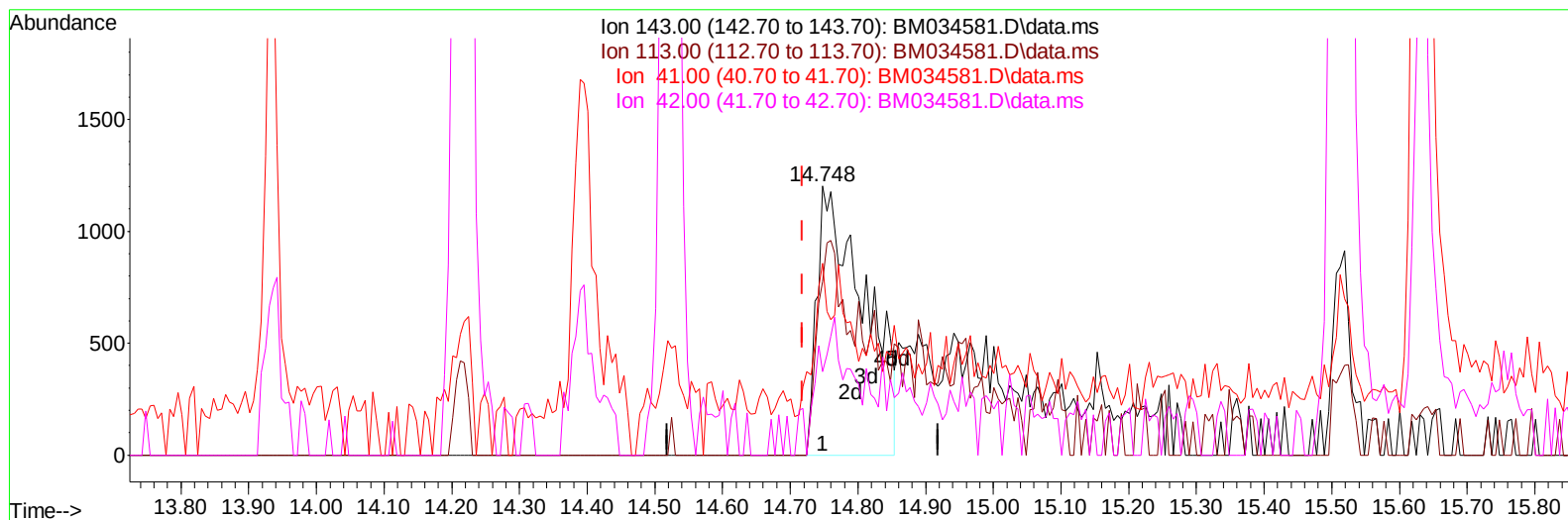
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(54) 4-Nitrophenol-d4 (S)

14.748min (+ 0.030) 1.67 ng/ul m

response 5800

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	77.80	64.48
41.00	31.90	71.30#
42.00	19.10	31.03#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.878	152	137851	20.000	ng/ul	0.00
20) Naphthalene-d8	10.684	136	550245	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.524	164	314905	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.265	188	612408	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.442	240	544775	20.000	ng/ul	# 0.00
88) Perylene-d12	23.818	264	591690	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.255	96	15201	3.932	ng/uL	0.00
4) Pyridine-d5	3.684	84	58259	6.245	ng/ul	0.00
7) Phenol-d5	7.037	99	60350	5.260	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.201	67	240676	31.479	ng/ul	0.00
11) 2-Chlorophenol-d4	7.407	132	210804	23.767	ng/ul	0.00
15) 4-Methylphenol-d8	8.590	113	121704	13.579	ng/ul	0.00
21) Nitrobenzene-d5	9.037	128	132539	32.774	ng/ul	0.00
24) 2-Nitrophenol-d4	9.766	143	135851	32.266	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.307	165	211731	26.707	ng/ul	0.00
31) 4-Chloroaniline-d4	10.825	131	248793	22.747	ng/ul	0.00
46) Dimethylphthalate-d6	13.936	166	778350	34.207	ng/ul	0.00
49) Acenaphthylene-d8	14.219	160	936491	32.021	ng/ul	0.00
54) 4-Nitrophenol-d4	14.748	143	5800m	1.672	ng/ul	0.03
60) Fluorene-d10	15.513	176	675384	34.347	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.636	200	103569	27.703	ng/ul	0.00
73) Anthracene-d10	17.365	188	1067766	36.570	ng/ul	0.00
81) Pyrene-d10	19.654	212	1284097	43.187	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.665	264	1118087	37.566	ng/ul	0.00

Target Compounds Qvalue

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

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