

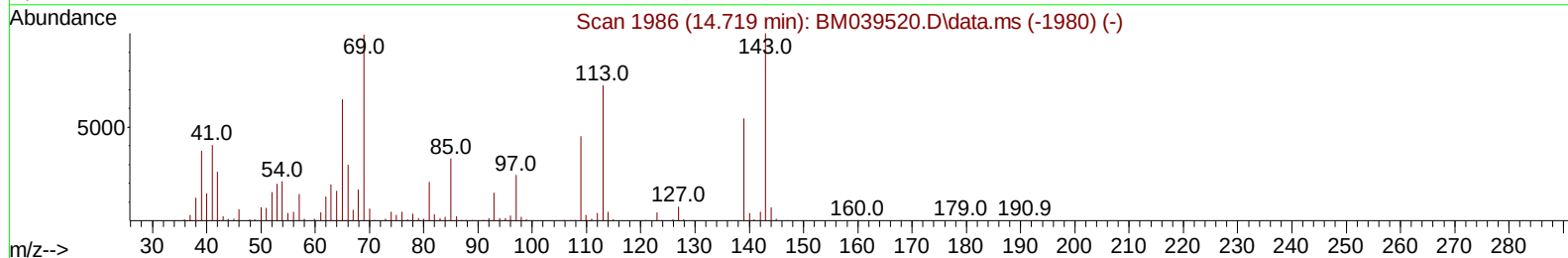
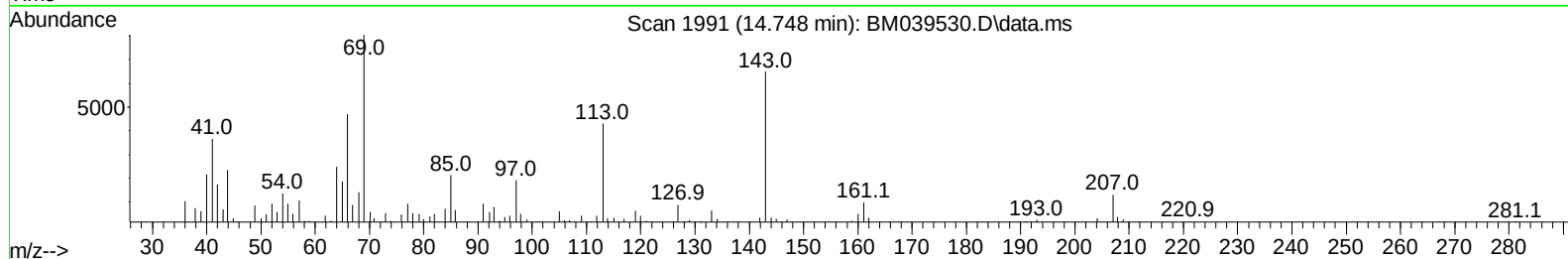
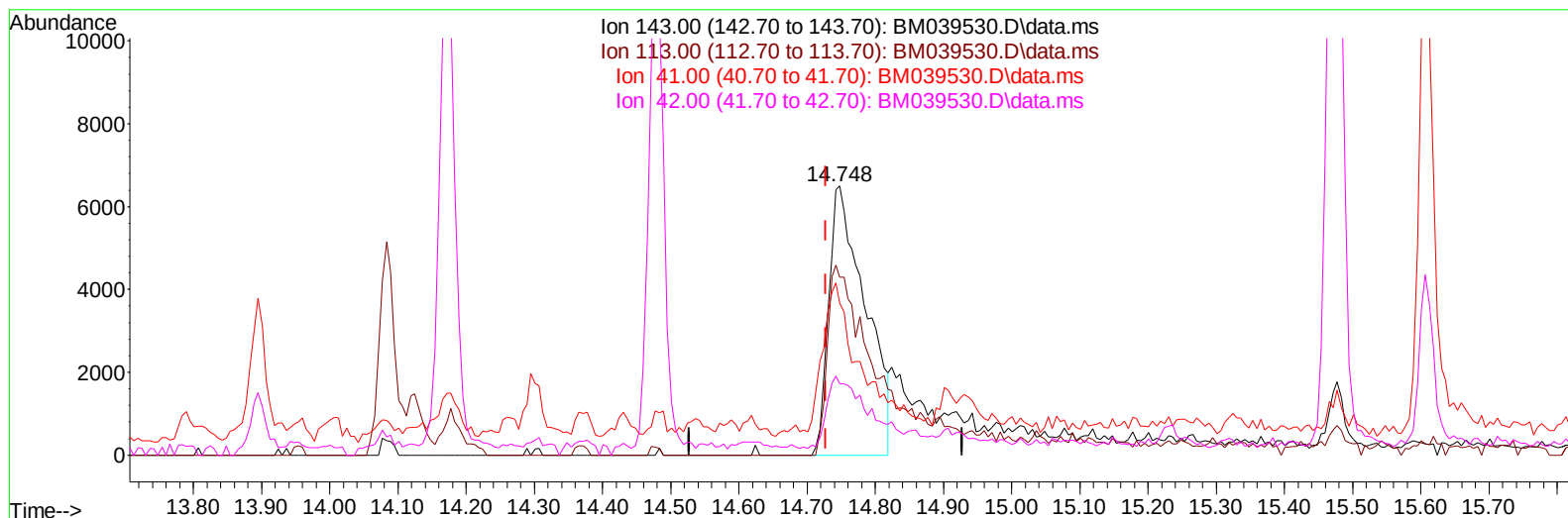
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042023\
Data File : BM039530.D
Acq On : 20 Apr 2023 18:25
Operator : CG/JU
Sample : 02378-07
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COMK0

Manual IntegrationsAPPROVED

Quant Time: Apr 21 00:38:47 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 19 08:18:41 2023
Response via : Initial Calibration

Reviewed By :Christian Giraldo 04/21/2023
Supervised By :Jagrut Upadhyay 04/21/2023



TIC: BM039530.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.748min (+ 0.021) 3.14 ng/ul

response 24253

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	66.12
41.00	39.90	56.58#
42.00	26.60	26.55

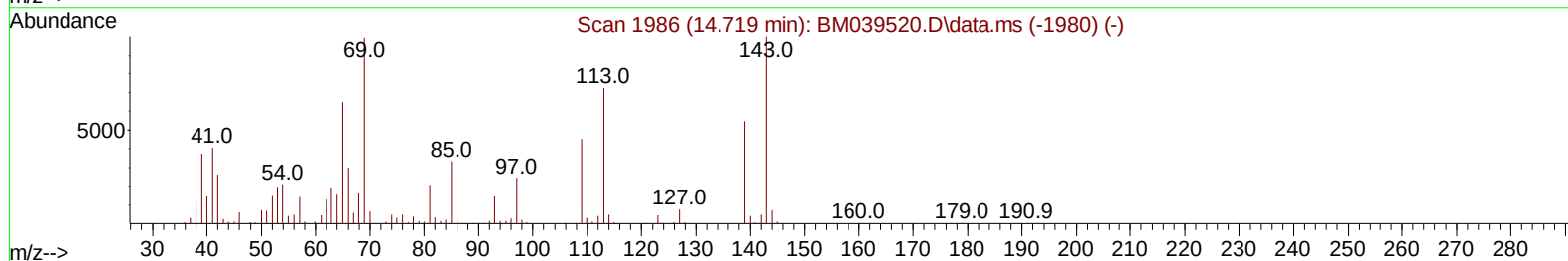
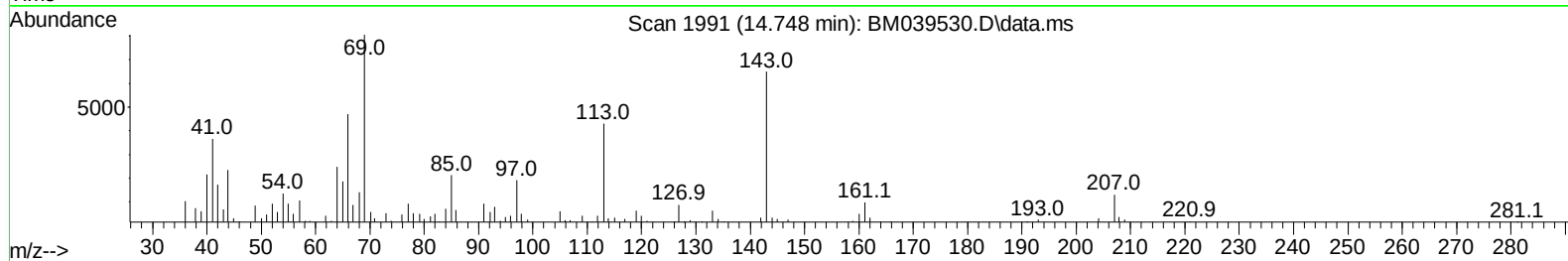
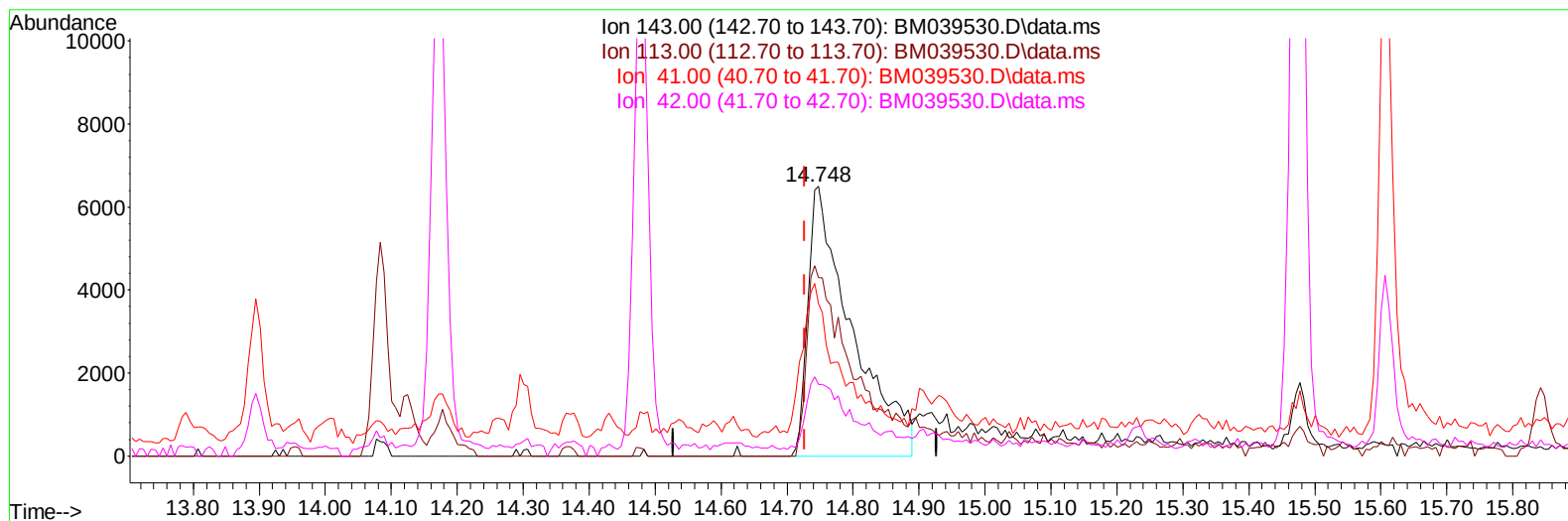
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(54) 4-Nitrophenol-d4 (S)

14.748min (+ 0.021) 3.90 ng/ul m

response 30123

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	66.12
41.00	39.90	56.58#
42.00	26.60	26.55

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.819	152	218180	20.000	ng/ul	0.00
20) Naphthalene-d8	10.631	136	985398	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.477	164	632512	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.230	188	1251528	20.000	ng/ul	0.00
79) Chrysene-d12	21.424	240	843586	20.000	ng/ul	0.00
88) Perylene-d12	23.788	264	867059	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	25543	4.494	ng/uL	0.00
4) Pyridine-d5	3.654	84	71373	4.469	ng/ul	0.01
7) Phenol-d5	7.013	99	119448	5.982	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.154	67	353468	26.780	ng/ul	0.00
11) 2-Chlorophenol-d4	7.360	132	362940	23.170	ng/ul	0.00
15) 4-Methylphenol-d8	8.554	113	240154	14.930	ng/ul	0.00
21) Nitrobenzene-d5	8.995	128	225410	28.433	ng/ul	0.00
24) 2-Nitrophenol-d4	9.719	143	255184	28.666	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.266	165	402761	26.294	ng/ul	0.00
31) 4-Chloroaniline-d4	10.795	131	324411	14.181	ng/ul	0.00
46) Dimethylphthalate-d6	13.895	166	1599584	31.357	ng/ul	0.00
49) Acenaphthylene-d8	14.171	160	1655037	28.880	ng/ul	0.00
54) 4-Nitrophenol-d4	14.748	143	30123m	3.902	ng/ul	0.02
60) Fluorene-d10	15.477	176	1286940	30.793	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.607	200	216453	26.026	ng/ul	0.00
73) Anthracene-d10	17.330	188	1902899	31.616	ng/ul	0.00
81) Pyrene-d10	19.630	212	1874138	33.433	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.636	264	1528727	32.657	ng/ul	0.00

Target Compounds Qvalue

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

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