

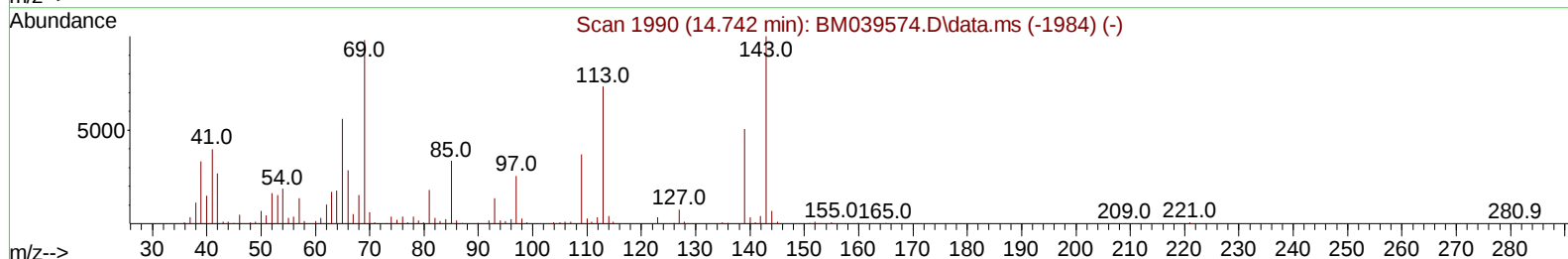
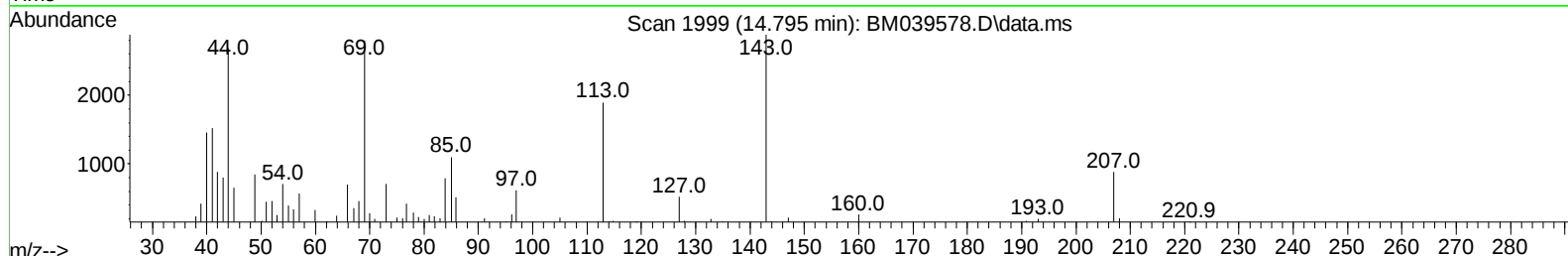
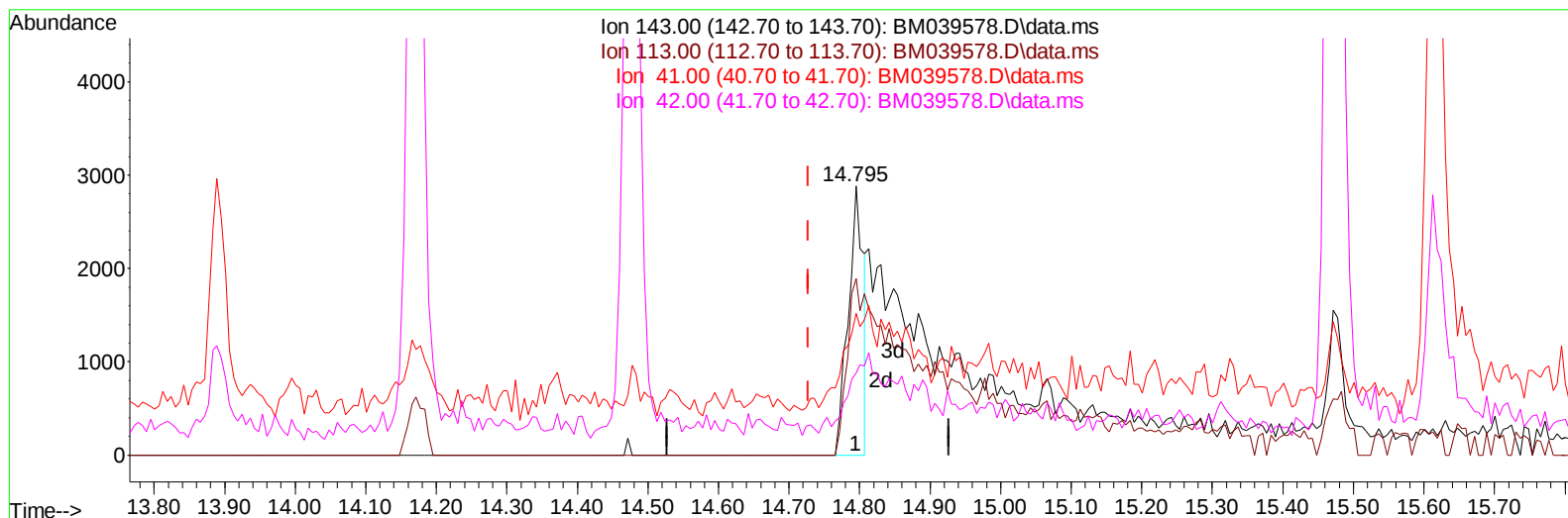
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042223\
Data File : BM039578.D
Acq On : 22 Apr 2023 16:16
Operator : CG/JU
Sample : 02440-03
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COMP8

Manual IntegrationsAPPROVED

Quant Time: Apr 22 23:34:17 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 04/24/2023
Supervised By :Jagrut Upadhyay 04/24/2023



TIC: BM039578.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.795min (+ 0.068) 0.84 ng/u1

response 4268

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	65.75
41.00	39.90	52.67#
42.00	26.60	30.57

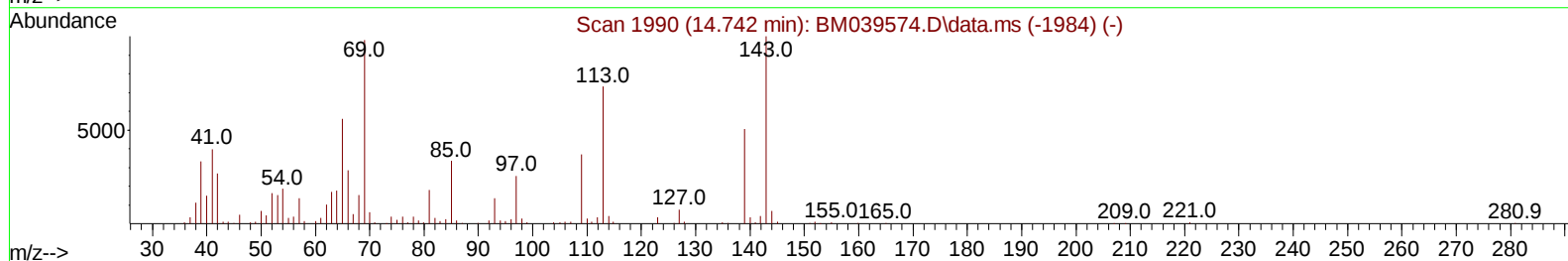
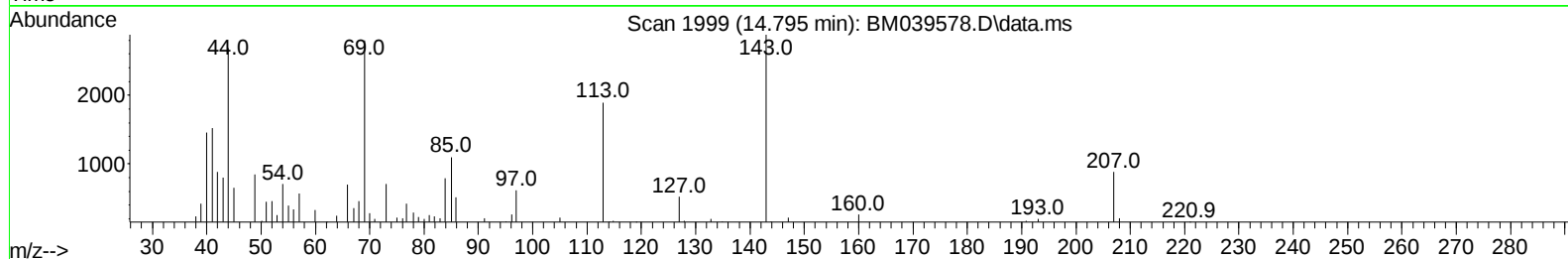
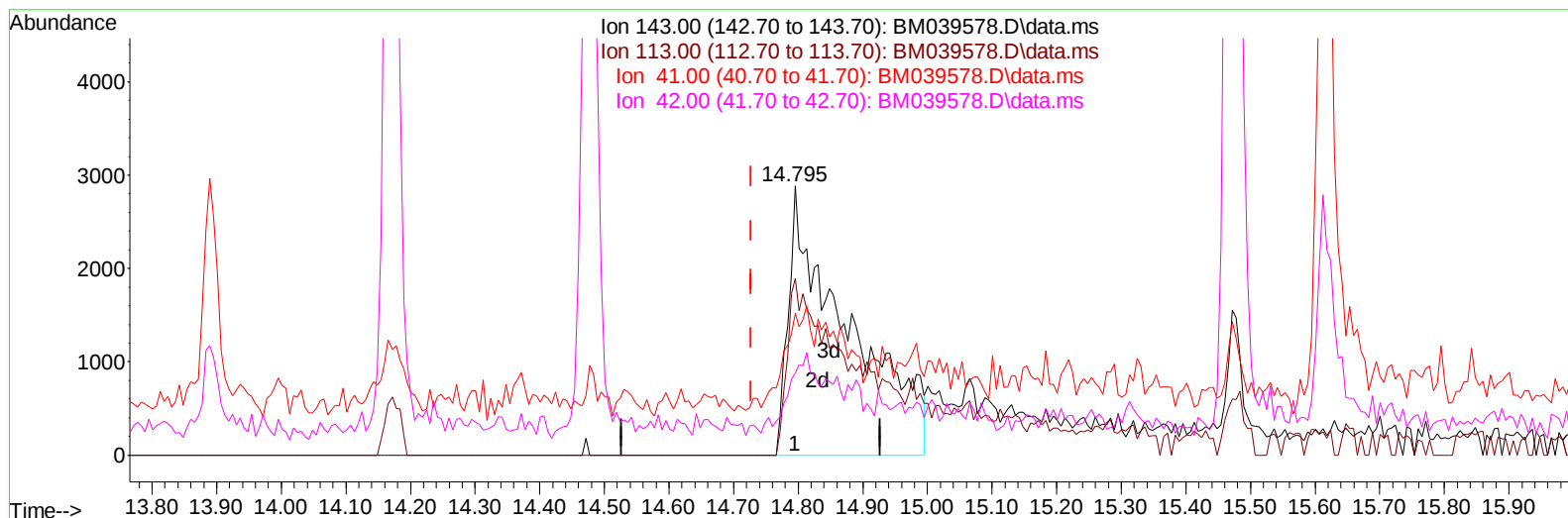
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(54) 4-Nitrophenol-d4 (S)

14.795min (+ 0.068) 3.60 ng/ul m

response 18251

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	65.75
41.00	39.90	52.67#
42.00	26.60	30.57

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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	7.819	152	175980	20.000	ng/ul	0.00
20) Naphthalene-d8	10.625	136	719190	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.477	164	415778	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.230	188	821328	20.000	ng/ul	0.00
79) Chrysene-d12	21.424	240	703786	20.000	ng/ul	0.00
88) Perylene-d12	23.794	264	771690	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	21637	4.719	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.013	99	94648	5.876	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.154	67	302570	28.421	ng/ul	0.00
11) 2-Chlorophenol-d4	7.354	132	302276	23.924	ng/ul	0.00
15) 4-Methylphenol-d8	8.554	113	170068	13.108	ng/ul	0.00
21) Nitrobenzene-d5	8.989	128	181121	31.303	ng/ul	0.00
24) 2-Nitrophenol-d4	9.713	143	197647	30.421	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.272	165	288528	25.809	ng/ul	0.00
31) 4-Chloroaniline-d4	10.789	131	96496	5.779	ng/ul	0.00
46) Dimethylphthalate-d6	13.895	166	1131808	33.753	ng/ul	0.00
49) Acenaphthylene-d8	14.172	160	1238341	32.873	ng/ul	0.00
54) 4-Nitrophenol-d4	14.795	143	18251m	3.596	ng/ul	0.07
60) Fluorene-d10	15.471	176	944963	34.397	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.618	200	135981	24.914	ng/ul	0.00
73) Anthracene-d10	17.330	188	1518660	38.448	ng/ul	0.00
81) Pyrene-d10	19.630	212	1775069	37.956	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.641	264	1624164	38.984	ng/ul	0.00

Target Compounds Qvalue

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

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