

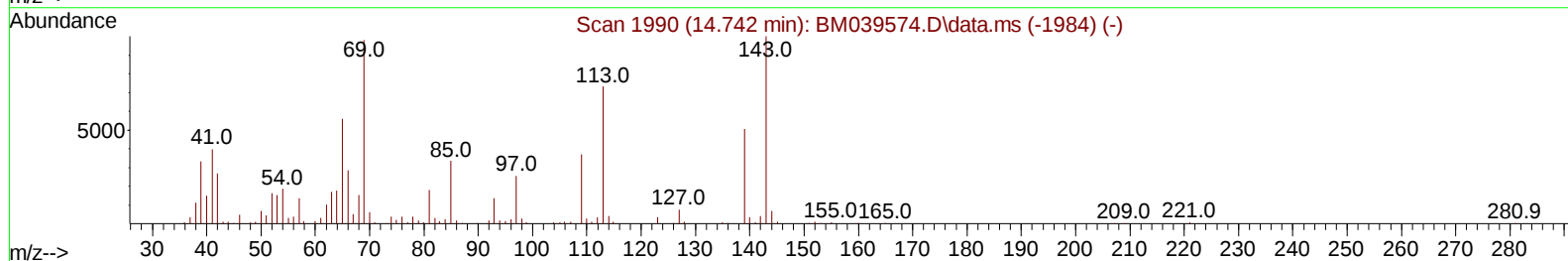
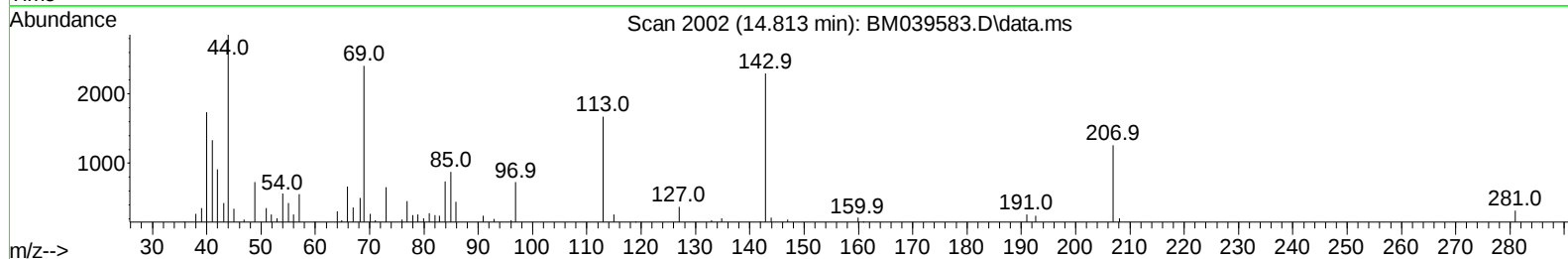
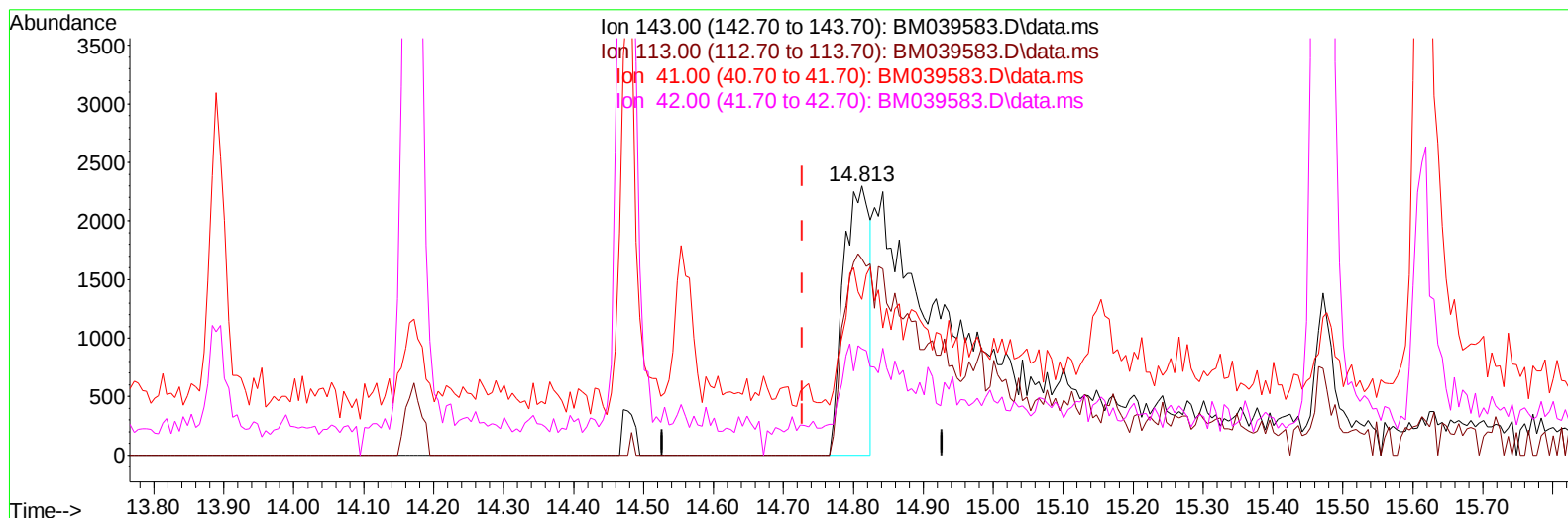
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042223\
Data File : BM039583.D
Acq On : 22 Apr 2023 19:17
Operator : CG/JU
Sample : 02440-13
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
COMQ8

Quant Time: Apr 22 23:34:58 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

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/24/2023
Supervised By :Jagrut Upadhyay 04/24/2023



TIC: BM039583.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.813min (+ 0.085) 1.16 ng/u1

response 6027

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	72.93
41.00	39.90	57.96#
42.00	26.60	39.60#

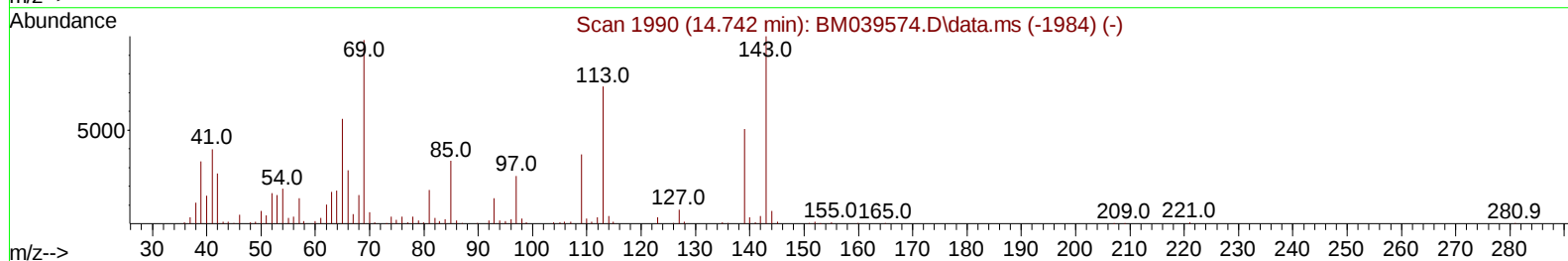
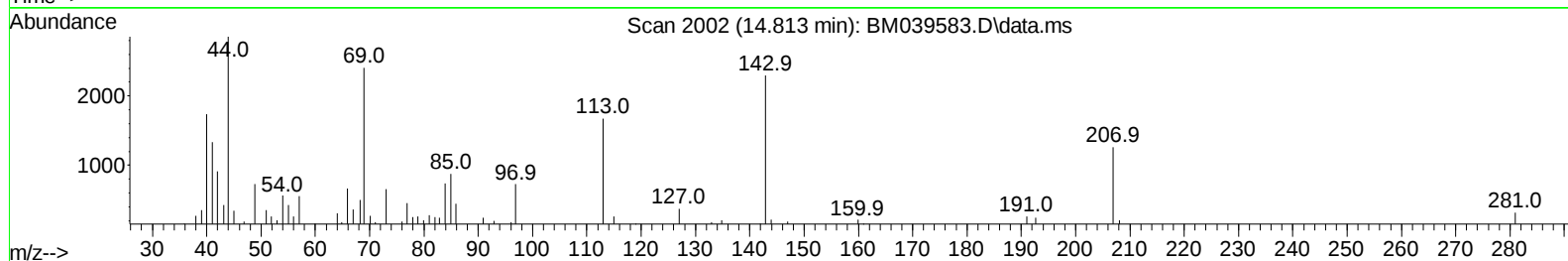
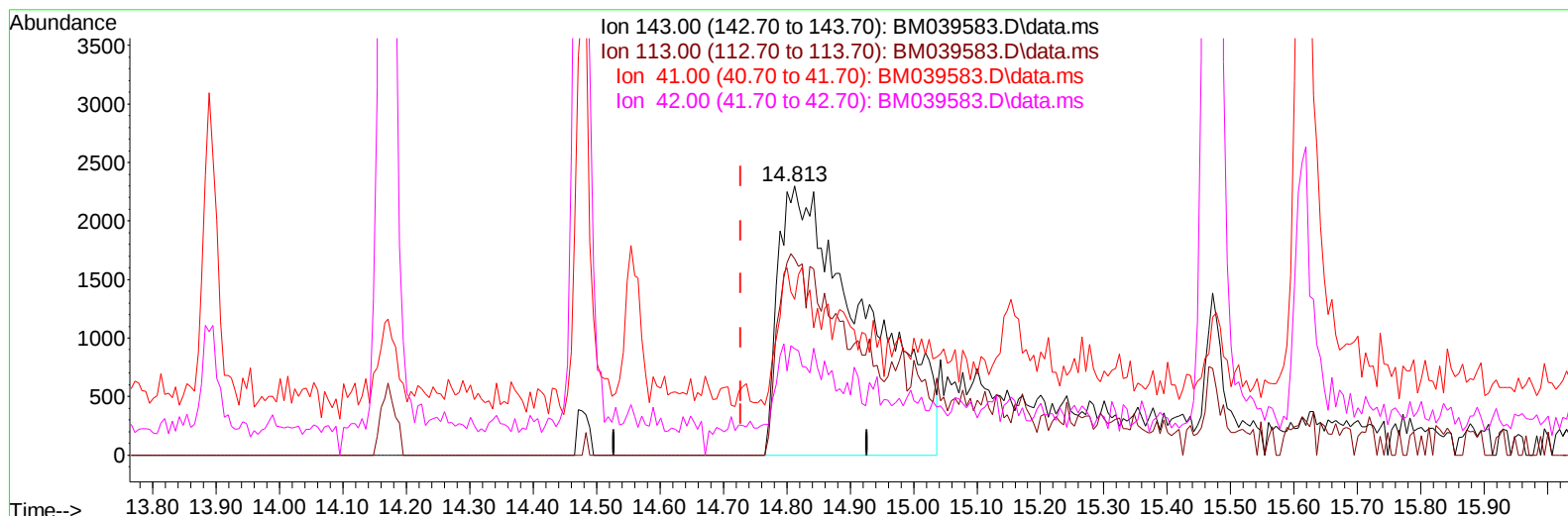
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(54) 4-Nitrophenol-d4 (S)

14.813min (+ 0.085) 4.16 ng/ul m

response 21643

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	72.93
41.00	39.90	57.96#
42.00	26.60	39.60#

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.819	152	185489	20.000	ng/ul	0.00
20) Naphthalene-d8	10.625	136	738658	20.000	ng/ul	-0.01
38) Acenaphthene-d10	14.477	164	425776	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.230	188	826955	20.000	ng/ul	0.00
79) Chrysene-d12	21.424	240	707331	20.000	ng/ul	0.00
88) Perylene-d12	23.788	264	781003	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.213	96	25498	5.276	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.013	99	111297	6.556	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.154	67	339018	30.213	ng/ul	0.00
11) 2-Chlorophenol-d4	7.354	132	339062	25.460	ng/ul	0.00
15) 4-Methylphenol-d8	8.554	113	205991	15.063	ng/ul	0.00
21) Nitrobenzene-d5	8.989	128	204381	34.392	ng/ul	0.00
24) 2-Nitrophenol-d4	9.713	143	223381	33.476	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.266	165	319758	27.849	ng/ul	0.00
31) 4-Chloroaniline-d4	10.783	131	151781	8.851	ng/ul	0.00
46) Dimethylphthalate-d6	13.889	166	1197542	34.875	ng/ul	-0.01
49) Acenaphthylene-d8	14.171	160	1360071	35.257	ng/ul	0.00
54) 4-Nitrophenol-d4	14.813	143	21643m	4.165	ng/ul	0.09
60) Fluorene-d10	15.471	176	1004903	35.719	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.613	200	144692	26.330	ng/ul	0.00
73) Anthracene-d10	17.330	188	1649539	41.477	ng/ul	0.00
81) Pyrene-d10	19.630	212	1864822	39.675	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.636	264	1743388	41.346	ng/ul	0.00

Target Compounds Qvalue

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

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