

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM042423\
Data File : BM039615.D
Acq On : 24 Apr 2023 14:09
Operator : CG/JU
Sample : PB152282BL
Misc :
ALS Vial : 9 Sample Multiplier: 1

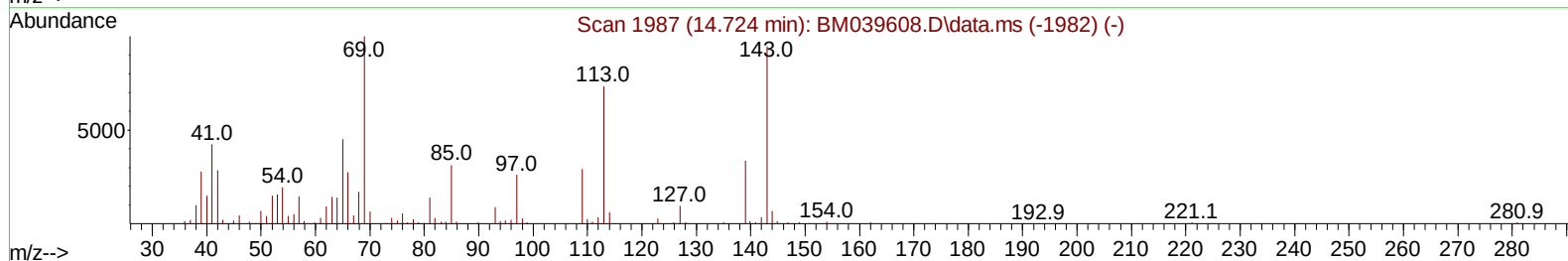
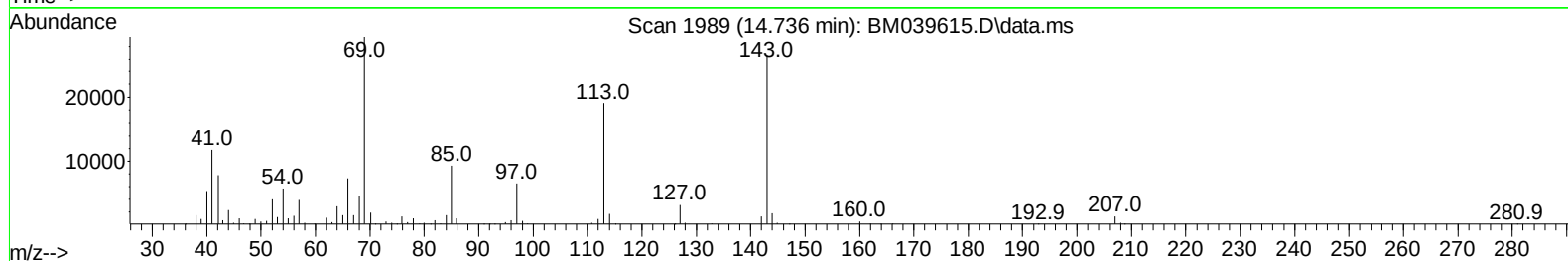
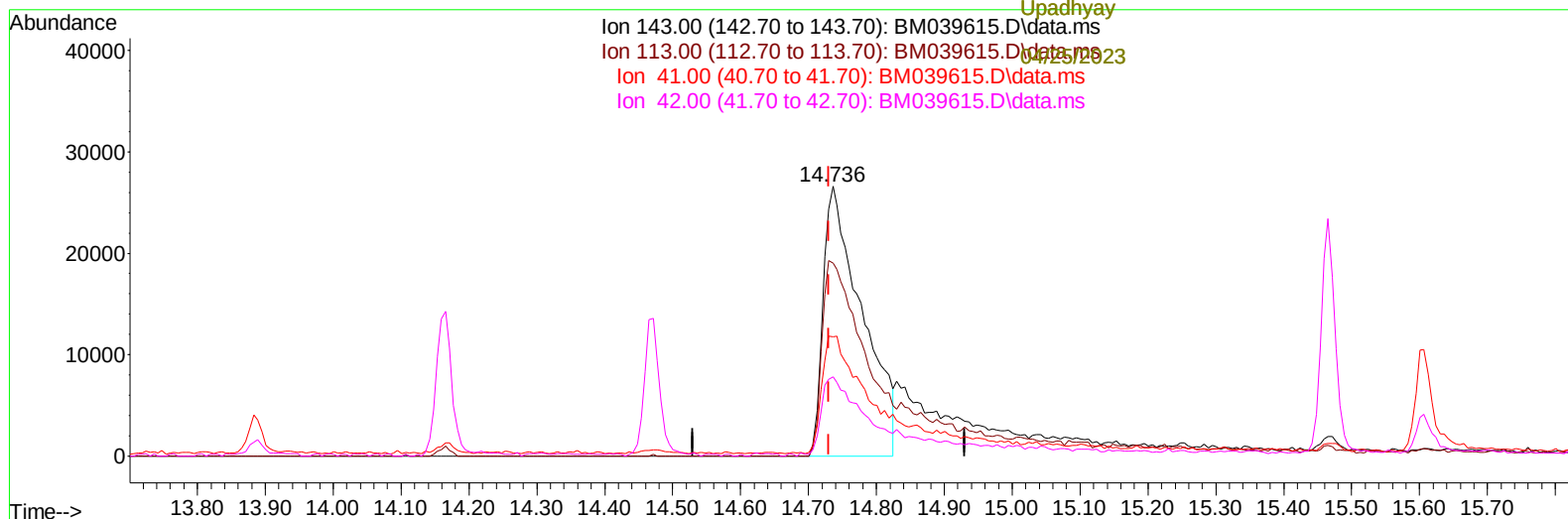
Instrument :
BNA_M
ClientSampleId :
SBLK282

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 04/25/2023
Supervised By :Jagrut Upadhyay 04/25/2023

Quant Time: Apr 25 00:13:16 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Apr 23 00:18:16 2023
Response via : Initial Calibration

04/25/2023
Supervised By :Jagrut
Upadhyay



TIC: BM039615.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.736min (+ 0.006) 14.65 ng/ul

response 105397

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	71.63
41.00	39.90	44.26
42.00	26.60	29.43

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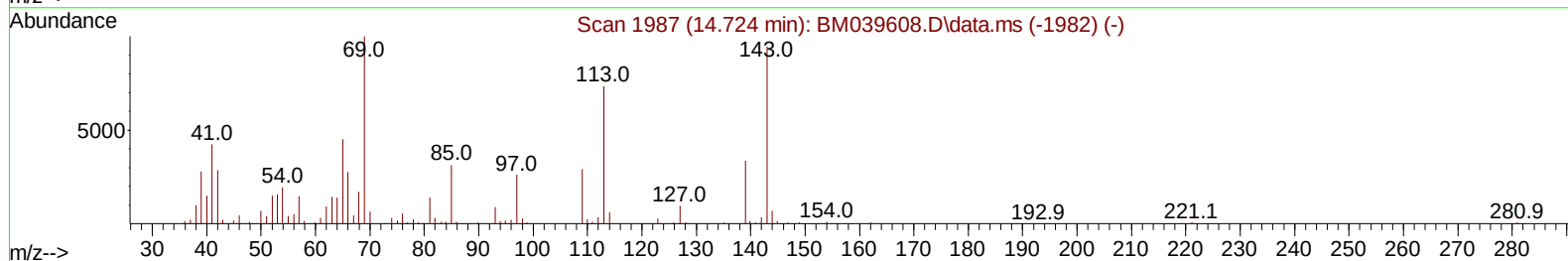
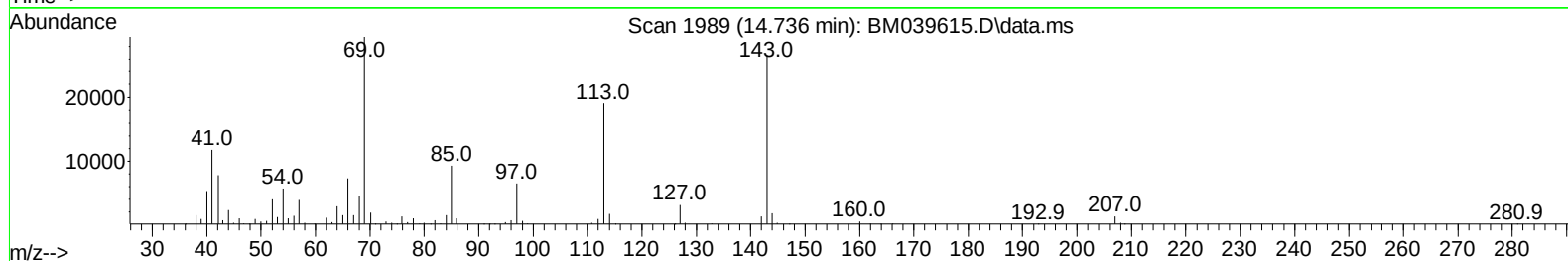
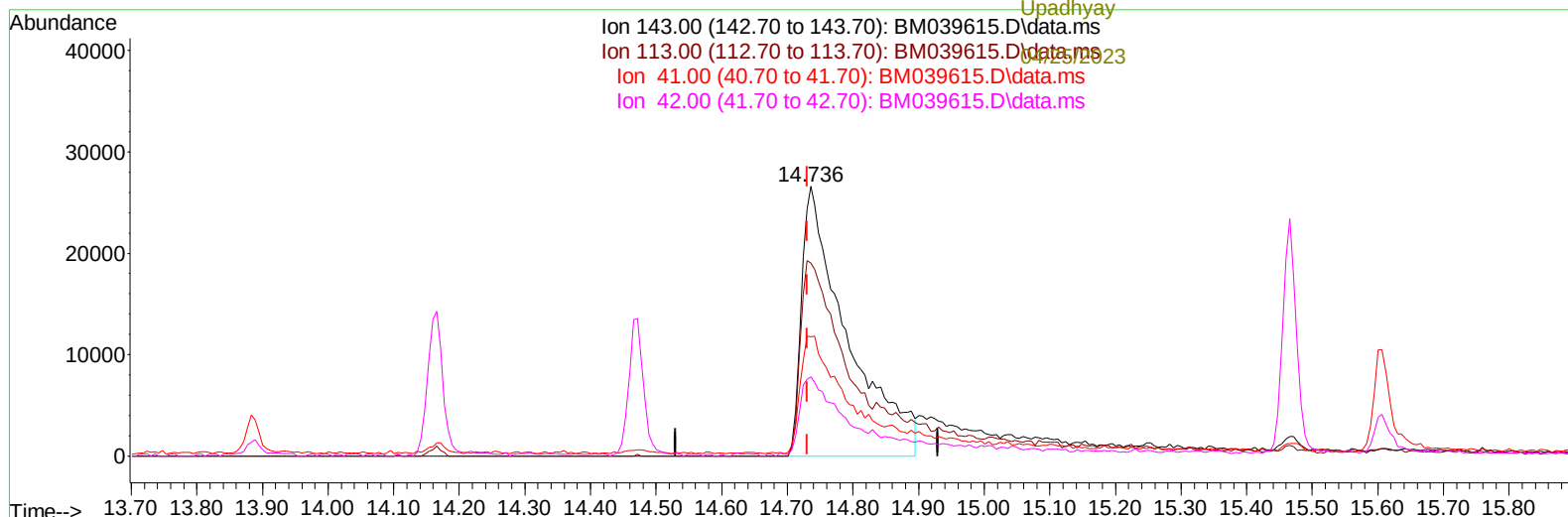
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TIC: BM039615.D\data.ms

(54) 4-Nitrophenol-d4 (S)

14.736min (+ 0.006) 17.76 ng/ul m

response 127740

Ion	Exp%	Act%
143.00	100.00	100.00
113.00	75.20	71.63
41.00	39.90	44.26
42.00	26.60	29.43

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Quant Time: Apr 25 00:38:59 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM041823.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sun Apr 23 00:18:16 2023
 Response via : Initial Calibration

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Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
-----04/25/2023						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.813	152	219273	20.000	ng/ul	0.00
20) Naphthalene-d8	10.619	136	948009	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.471	164	589248	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.224	188	1200008	20.000	ng/ul	0.00
79) Chrysene-d12	21.418	240	929804	20.000	ng/ul	0.00
88) Perylene-d12	23.782	264	940137	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.207	96	34795	6.091	ng/uL	0.00
4) Pyridine-d5	3.631	84	426871	26.594	ng/ul	0.00
7) Phenol-d5	6.995	99	534863	26.651	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.148	67	396534	29.894	ng/ul	0.00
11) 2-Chlorophenol-d4	7.342	132	451850	28.702	ng/ul	0.00
15) 4-Methylphenol-d8	8.542	113	437906	27.088	ng/ul	0.00
21) Nitrobenzene-d5	8.983	128	221510	29.043	ng/ul	0.00
24) 2-Nitrophenol-d4	9.707	143	250853	29.291	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.254	165	382934	25.986	ng/ul	0.00
31) 4-Chloroaniline-d4	10.772	131	615141	27.949	ng/ul	0.00
46) Dimethylphthalate-d6	13.889	166	1445651	30.421	ng/ul	0.00
49) Acenaphthylene-d8	14.165	160	1597841	29.929	ng/ul	0.00
54) 4-Nitrophenol-d4	14.736	143	127740m	17.761	ng/ul	0.00
60) Fluorene-d10	15.465	176	1195749	30.712	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.607	200	192004	24.077	ng/ul	0.00
73) Anthracene-d10	17.324	188	1867597	32.361	ng/ul	0.00
81) Pyrene-d10	19.624	212	2072642	33.546	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.630	264	1619122	31.899	ng/ul	0.00

Target Compounds Qvalue

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

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