

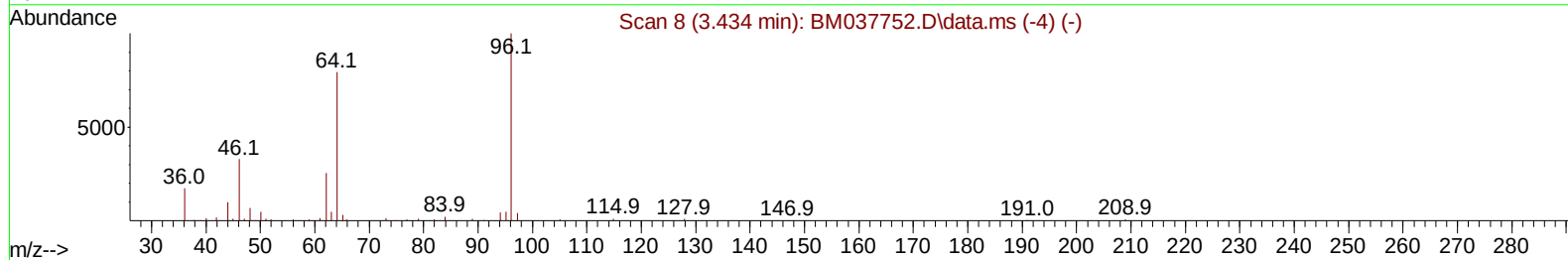
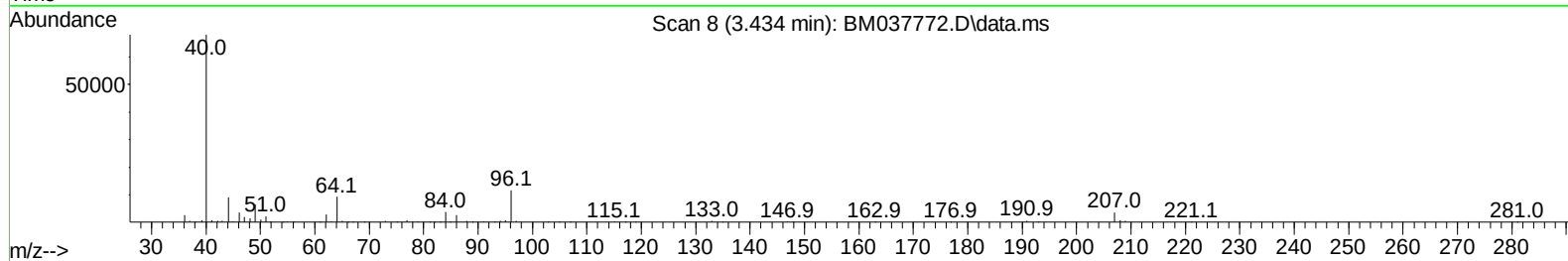
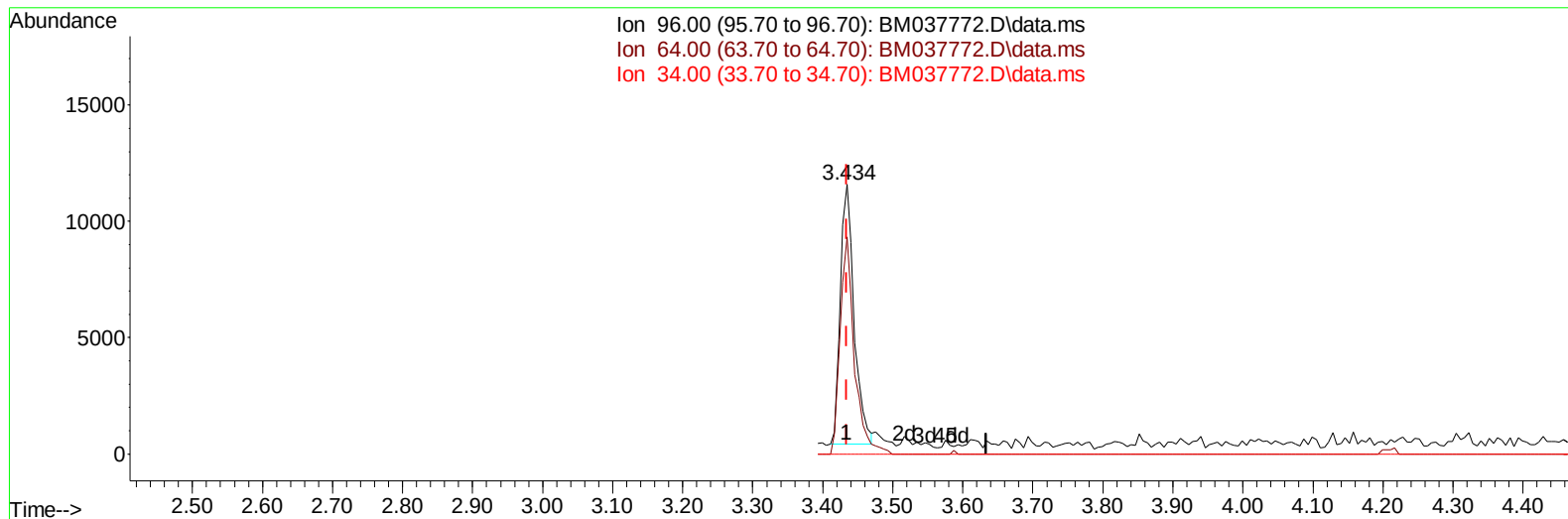
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM112822\
Data File : BM037772.D
Acq On : 28 Nov 2022 23:24
Operator : CG/JU
Sample : N5658-10
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
ESQT0

Manual IntegrationsAPPROVED

Quant Time: Nov 29 00:59:19 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM112322.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Nov 28 22:24:47 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/29/2022
Supervised By :mohammad ahmed 11/30/2022



TIC: BM037772.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.434min (+ 0.000) 3.79 ng/uL

response 15461

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	82.30	80.58
34.00	0.00	0.00
0.00	0.00	0.00

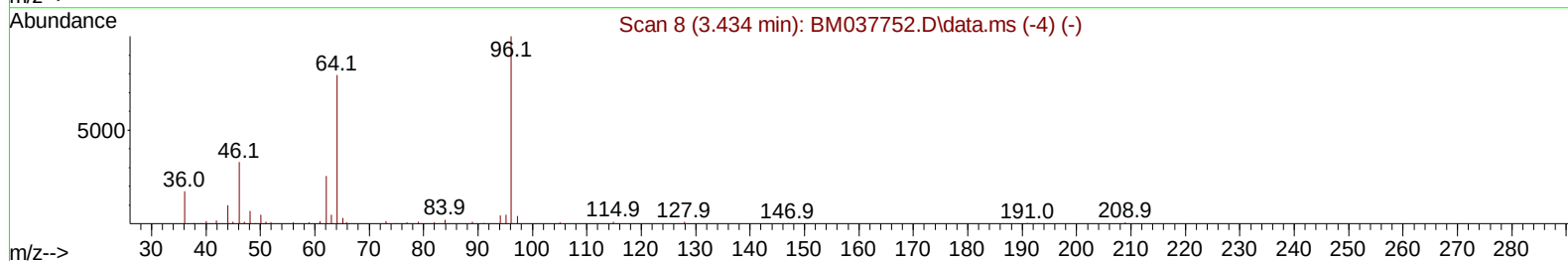
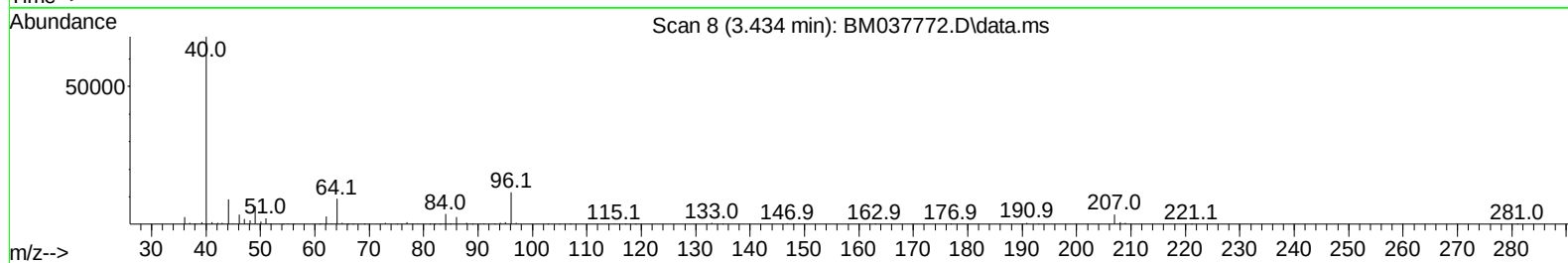
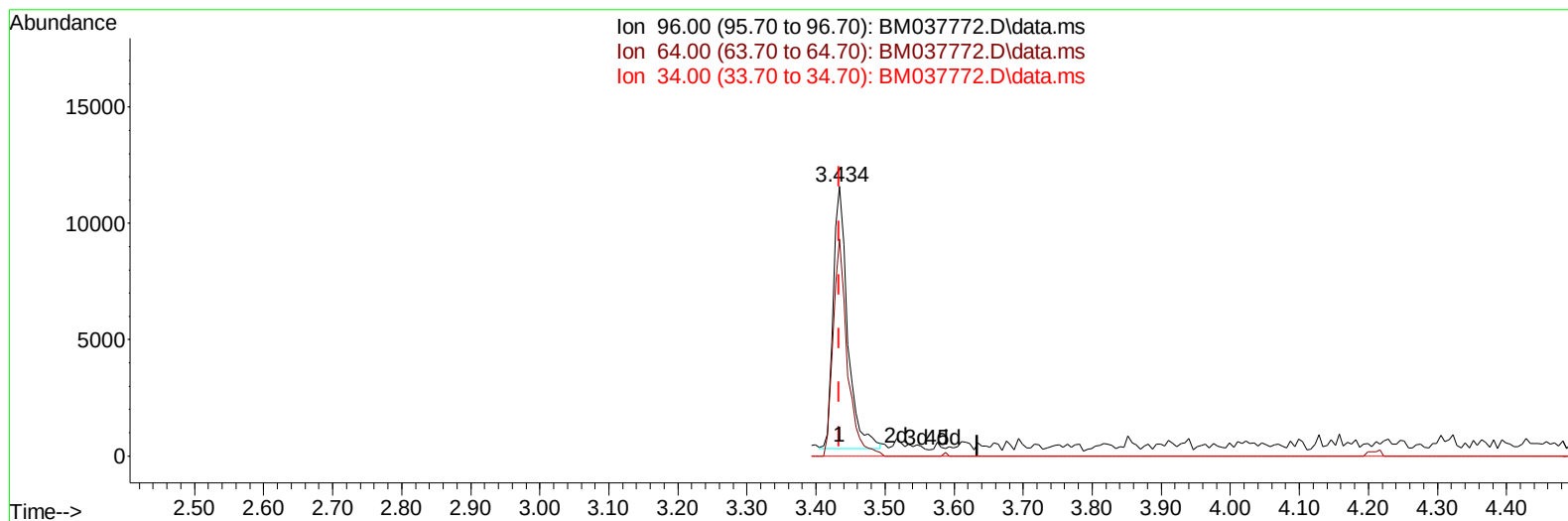
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TIC: BM037772.D\data.ms

(3) 1,4-Dioxane-d8 (S)

3.434min (+ 0.000) 4.03 ng/uL m

response 16451

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	82.30	80.58
34.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.105	152	169362	20.000	ng/ul	0.00
20) Naphthalene-d8	10.928	136	816164	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.734	164	503632	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.469	188	1010216	20.000	ng/ul	0.00
79) Chrysene-d12	21.633	240	970554	20.000	ng/ul	0.00
88) Perylene-d12	24.121	264	859492	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.434	96	16451m	4.028	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.252	99	336321	19.905	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.422	67	261567	23.207	ng/ul	0.00
11) 2-Chlorophenol-d4	7.628	132	267755	22.806	ng/ul	0.00
15) 4-Methylphenol-d8	8.810	113	226420	16.840	ng/ul	0.00
21) Nitrobenzene-d5	9.275	128	152897	23.896	ng/ul	0.00
24) 2-Nitrophenol-d4	10.004	143	146421	23.224	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.540	165	236837	20.528	ng/ul	0.00
31) 4-Chloroaniline-d4	11.063	131	111110	5.804	ng/ul	0.00
46) Dimethylphthalate-d6	14.140	166	825752	23.148	ng/ul	0.00
49) Acenaphthylene-d8	14.428	160	981260	23.304	ng/ul	0.00
54) 4-Nitrophenol-d4	14.910	143	71083	9.046	ng/ul	0.00
60) Fluorene-d10	15.716	176	731841	23.424	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.828	200	76342	14.048	ng/ul	0.00
73) Anthracene-d10	17.569	188	1127585	23.869	ng/ul	0.00
81) Pyrene-d10	19.845	212	1290856	26.172	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.957	264	909105	20.549	ng/ul	0.00

Target Compounds Qvalue

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

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