

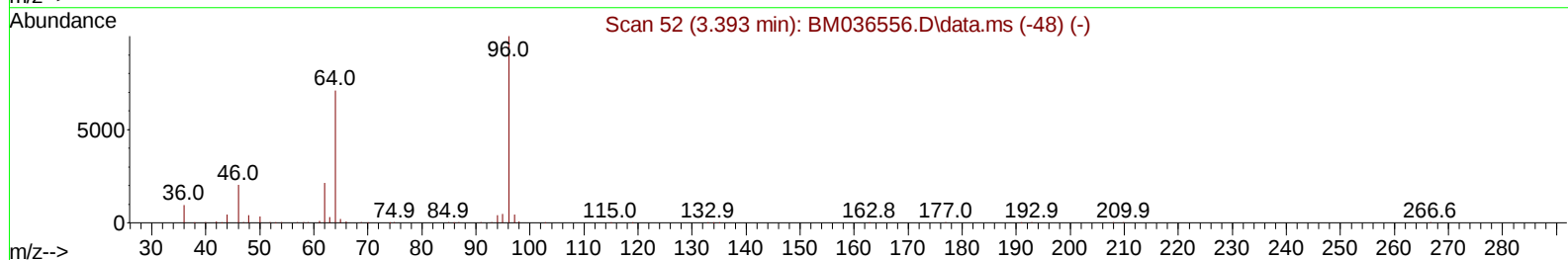
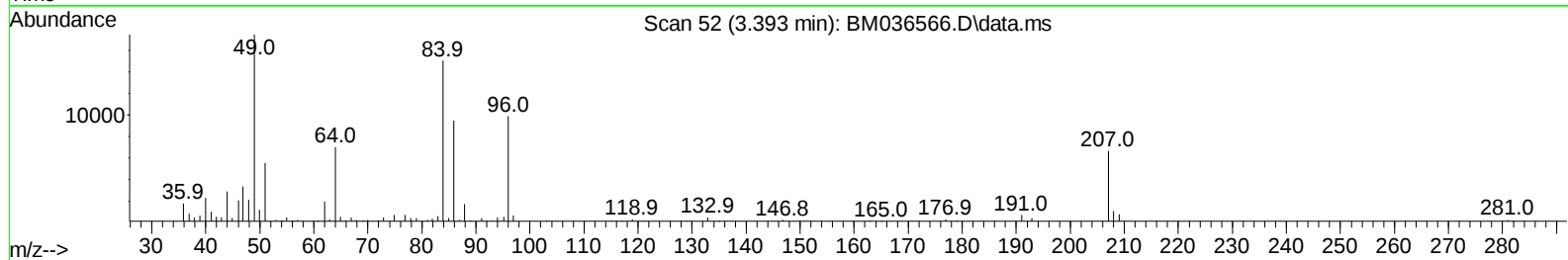
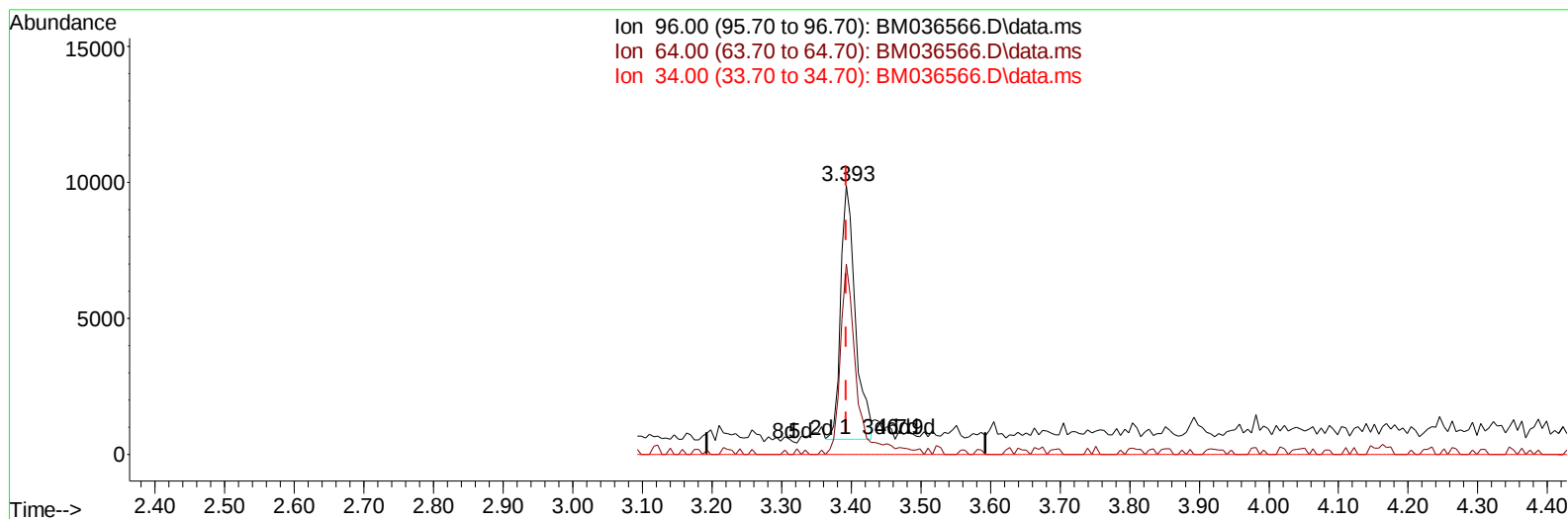
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM091622\
Data File : BM036566.D
Acq On : 16 Sep 2022 17:10
Operator : CG/JU
Sample : N4601-08
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A5749

Manual IntegrationsAPPROVED

Quant Time: Sep 16 23:01:07 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM091522.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Sep 16 02:39:54 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 09/19/2022
Supervised By :mohammad ahmed 09/19/2022



TIC: BM036566.D\data.ms

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

3.393min (-0.000) 1.43 ng/uL

response 13577

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.00	70.91
34.00	0.00	0.00
0.00	0.00	0.00

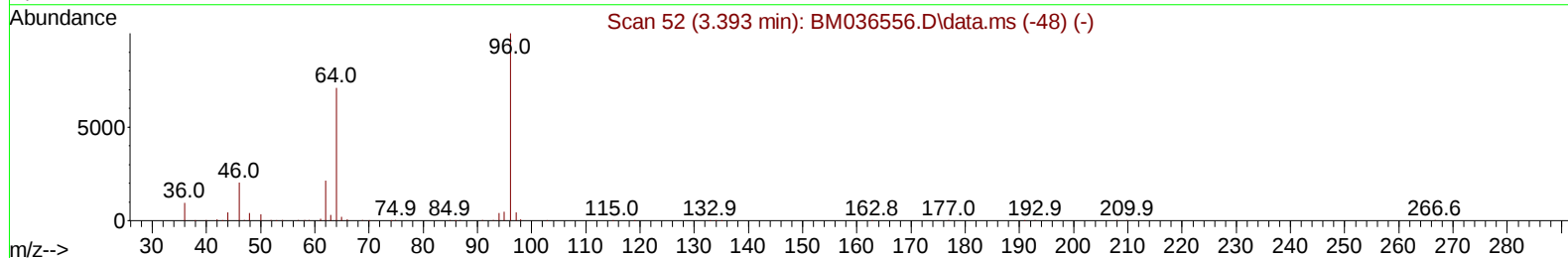
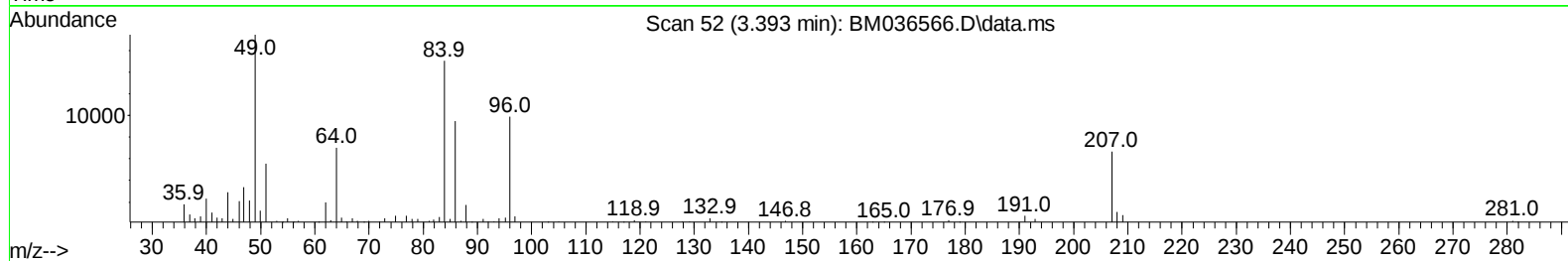
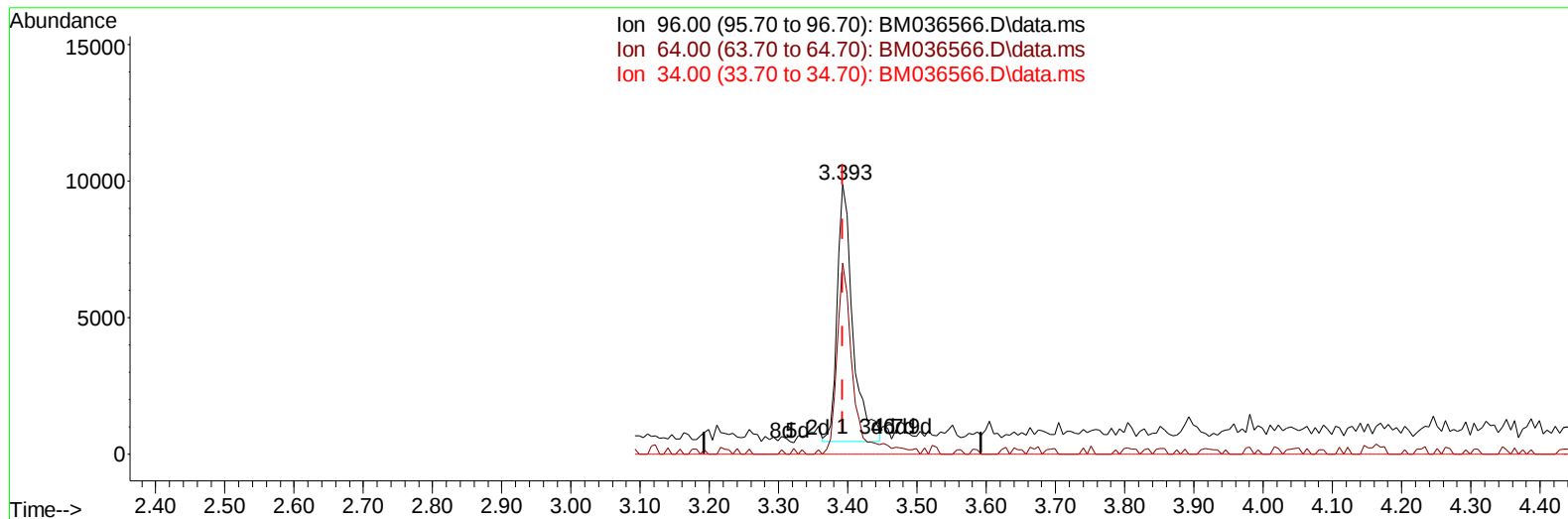
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TIC: BM036566.D\data.ms

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

3.393min (-0.000) 1.53 ng/uL m

response 14587

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.00	70.91
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	8.028	152	357129	20.000	ng/ul	0.00
20) Naphthalene-d8	10.839	136	1599169	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.651	164	1099200	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.386	188	2336163	20.000	ng/ul	0.00
79) Chrysene-d12	21.550	240	2303863	20.000	ng/ul	0.00
88) Perylene-d12	23.997	264	2328940	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.393	96	14587m	1.531	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/ul	
7) Phenol-d5	7.175	99	352360	11.041	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.351	67	258264	13.325	ng/ul	0.00
11) 2-Chlorophenol-d4	7.551	132	302475	13.164	ng/ul	0.00
15) 4-Methylphenol-d8	8.728	113	259546	11.006	ng/ul	0.00
21) Nitrobenzene-d5	9.192	128	148733	13.741	ng/ul	0.00
24) 2-Nitrophenol-d4	9.916	143	127428	13.143	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.451	165	273361	12.166	ng/ul	0.00
31) 4-Chloroaniline-d4	10.969	131	117709	3.151	ng/ul	0.00
46) Dimethylphthalate-d6	14.063	166	929609	12.539	ng/ul	0.00
49) Acenaphthylene-d8	14.345	160	1222474	13.902	ng/ul	0.00
54) 4-Nitrophenol-d4	14.821	143	52061	3.621	ng/ul	0.00
60) Fluorene-d10	15.639	176	980953	14.765	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.739	200	37898	4.100	ng/ul	-0.01
73) Anthracene-d10	17.486	188	1485829	13.910	ng/ul	0.00
81) Pyrene-d10	19.768	212	1690212	14.913	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.838	264	1572717	13.822	ng/ul	0.00

Target Compounds Qvalue

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

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