

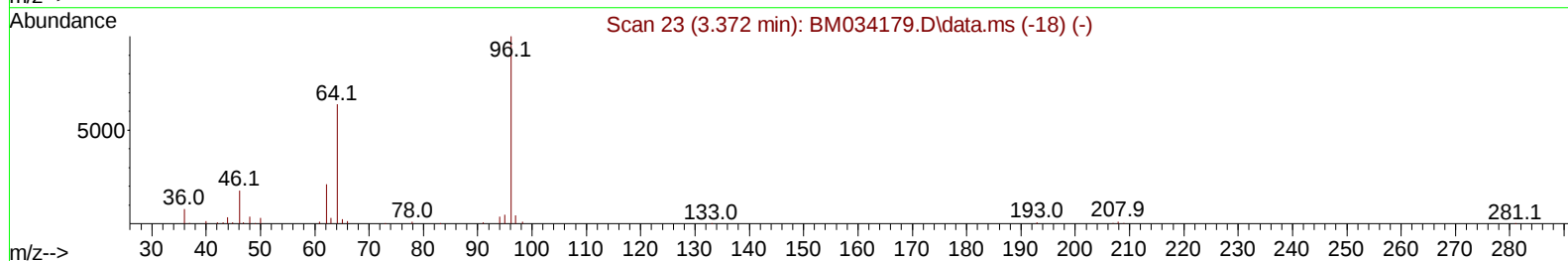
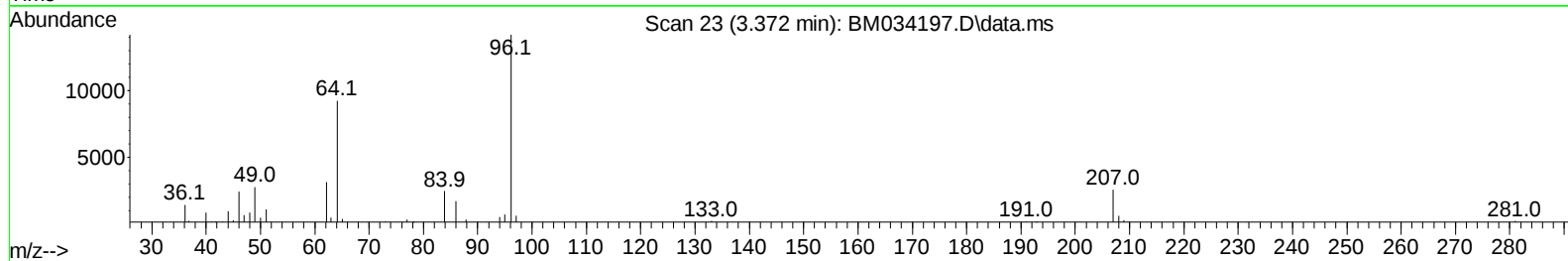
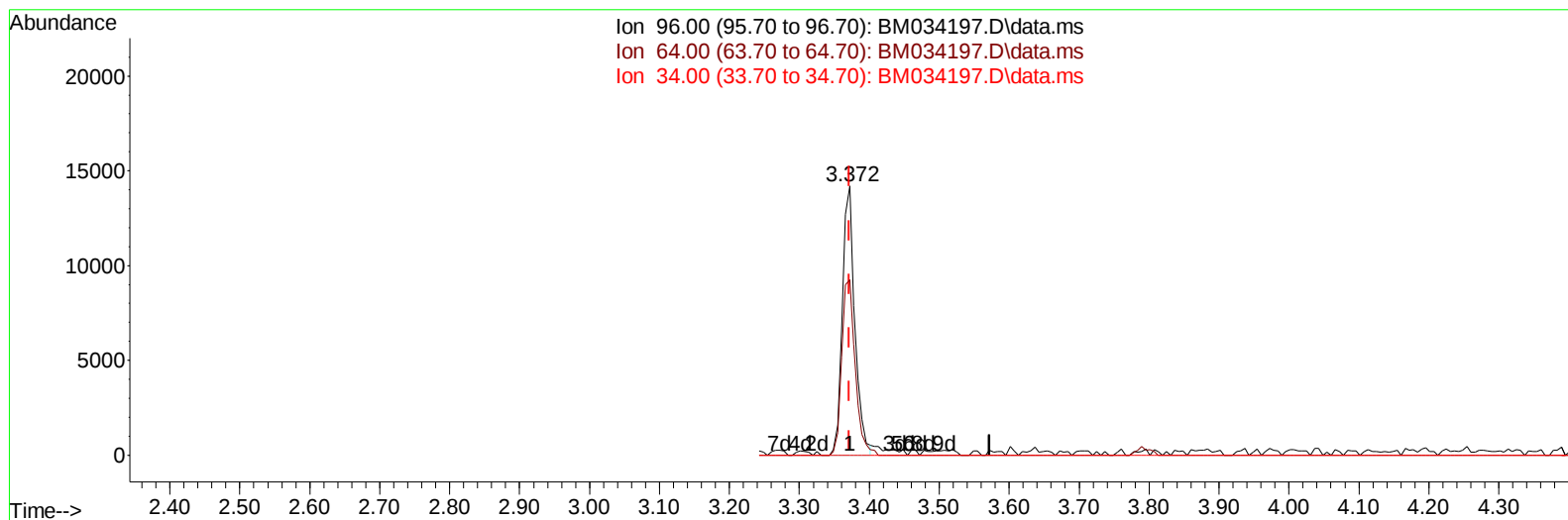
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
Data File : BM034197.D
Acq On : 09 Mar 2022 22:23
Operator : CG/JU
Sample : N1775-20
Misc :
ALS Vial : 57 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DBSC3

Manual IntegrationsAPPROVED

Quant Time: Mar 10 00:01:29 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Mar 08 13:09:16 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/10/2022
Supervised By :Yogesh Patel 03/11/2022



TIC: BM034197.D\data.ms

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

3.372min (+ 0.000) 5.59 ng/uL

response 17737

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	65.19
34.00	0.00	0.00
0.00	0.00	0.00

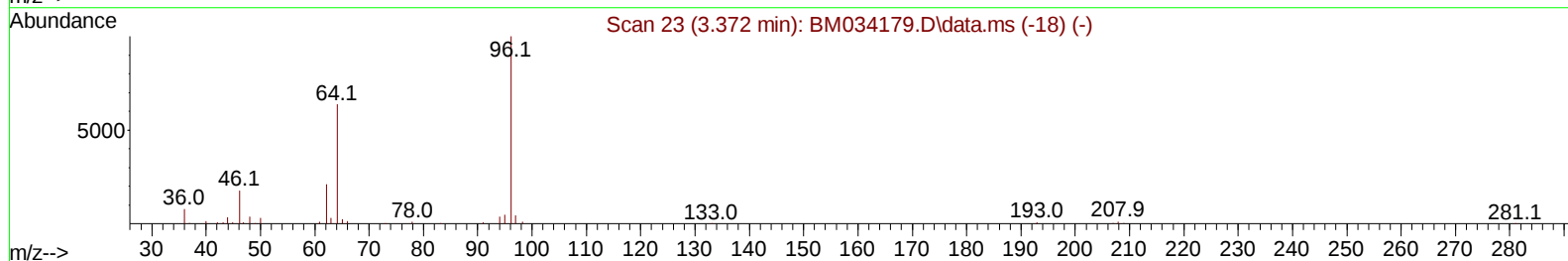
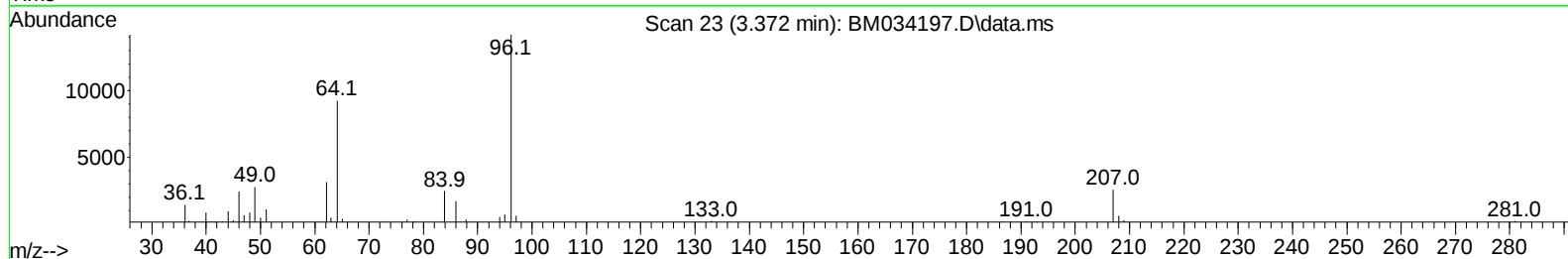
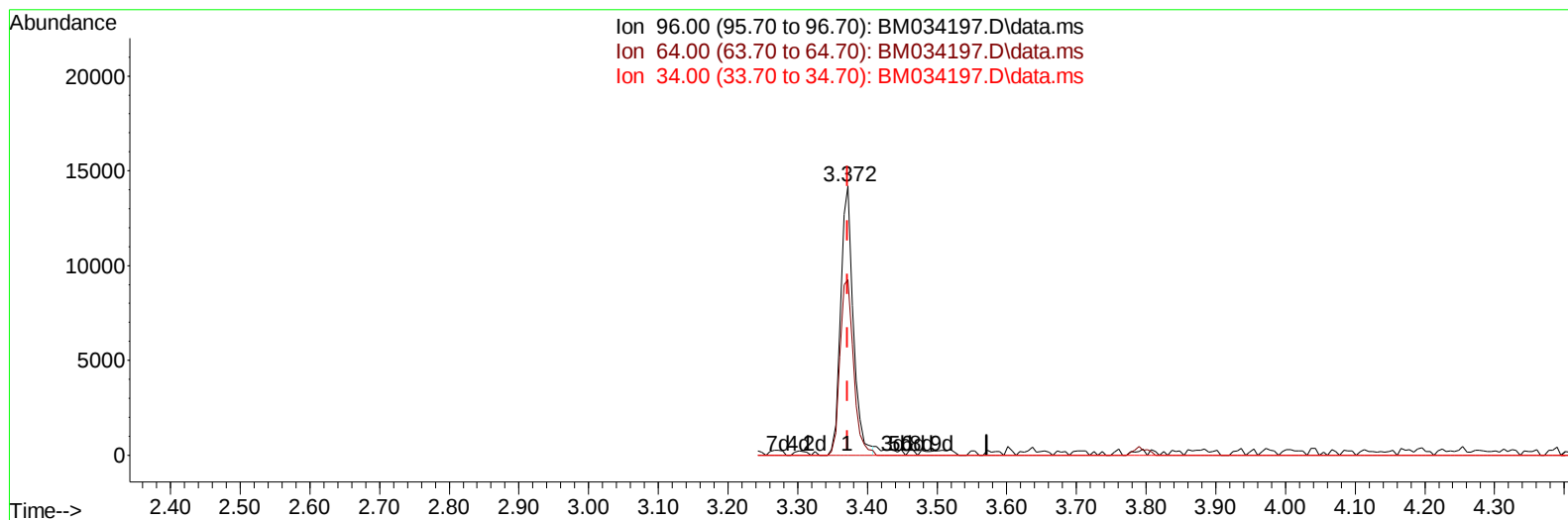
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TIC: BM034197.D\data.ms

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

3.372min (+ 0.000) 5.65 ng/uL m

response 17903

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	65.19
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	7.989	152	145837	20.000	ng/ul	0.00
20) Naphthalene-d8	10.801	136	587290	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.624	164	399661	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.365	188	783859	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.530	240	646908	20.000	ng/ul	-0.01
88) Perylene-d12	23.977	264	697914	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	17903m	5.646	ng/uL	0.00
4) Pyridine-d5	3.790	84	235651	25.530	ng/ul	0.00
7) Phenol-d5	7.137	99	321964	28.175	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.307	67	186280	30.287	ng/ul	0.00
11) 2-Chlorophenol-d4	7.513	132	284477	30.032	ng/ul	0.00
15) 4-Methylphenol-d8	8.689	113	280114	29.255	ng/ul	0.00
21) Nitrobenzene-d5	9.148	128	142688	31.357	ng/ul	0.00
24) 2-Nitrophenol-d4	9.878	143	158835	32.138	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.413	165	303969	29.824	ng/ul	0.00
31) 4-Chloroaniline-d4	10.930	131	396401	28.466	ng/ul	0.00
46) Dimethylphthalate-d6	14.030	166	999379	32.953	ng/ul	0.00
49) Acenaphthylene-d8	14.318	160	1206100	31.803	ng/ul	0.00
54) 4-Nitrophenol-d4	14.795	143	130291	23.089	ng/ul	-0.01
60) Fluorene-d10	15.612	176	853499	31.621	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.718	200	125976	24.904	ng/ul	0.00
73) Anthracene-d10	17.459	188	1255934	32.526	ng/ul	0.00
81) Pyrene-d10	19.748	212	1303129	36.647	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.818	264	1213500	32.777	ng/ul	-0.01

Target Compounds Qvalue

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

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