

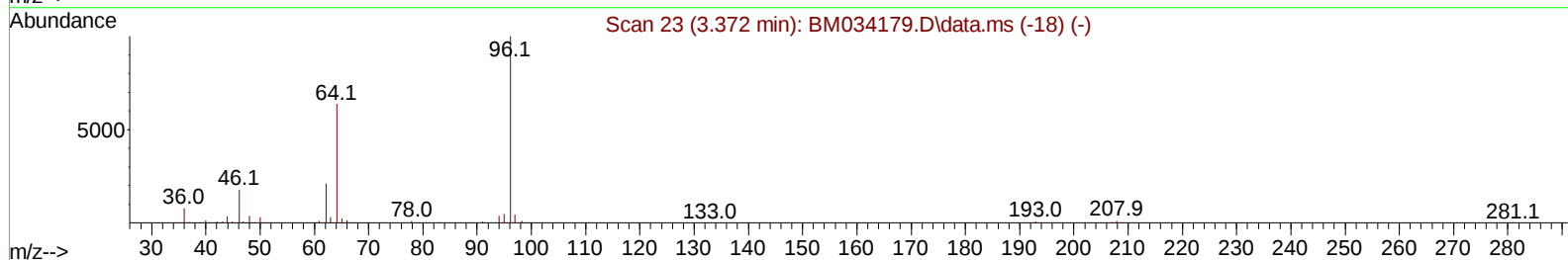
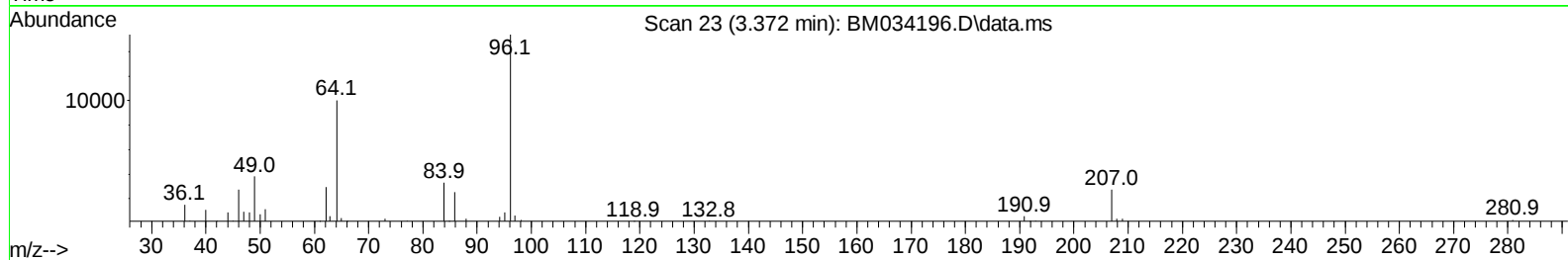
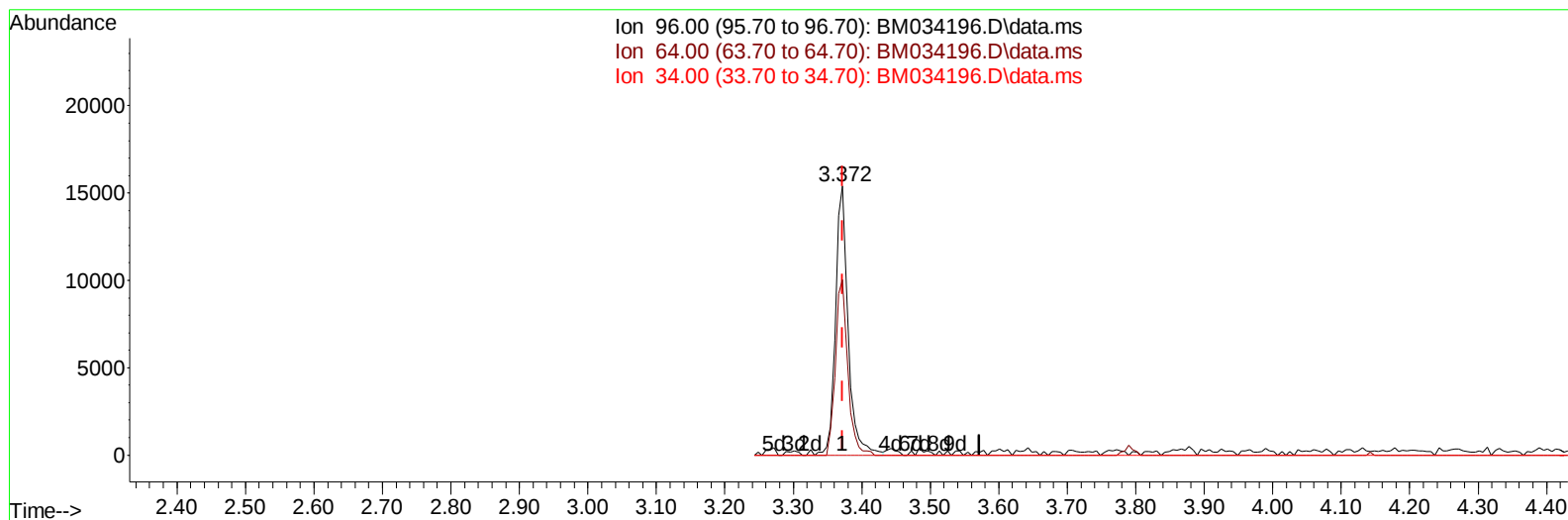
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034196.D
 Acq On : 09 Mar 2022 21:47
 Operator : CG/JU
 Sample : PB143147BL
 Misc :
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SBLK147

Manual IntegrationsAPPROVED

Quant Time: Mar 10 00:01:21 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: BM034196.D\data.ms

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

3.372min (+ 0.000) 5.86 ng/uL

response 19635

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	65.12
34.00	0.00	0.00
0.00	0.00	0.00

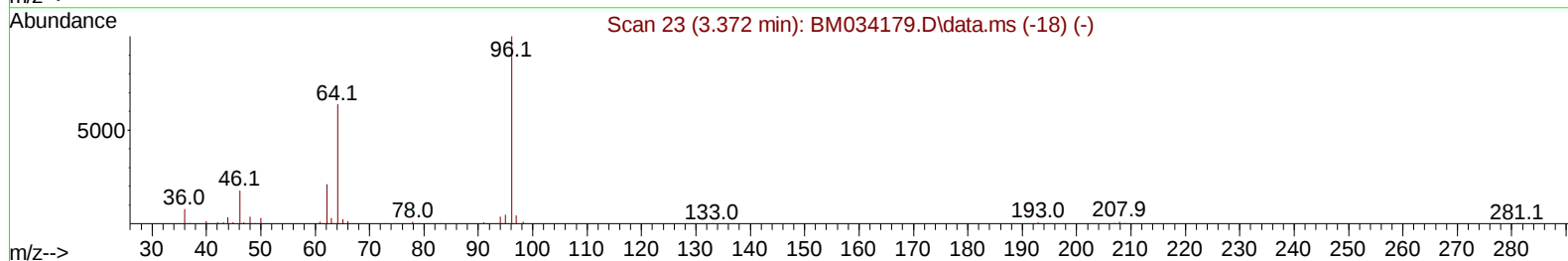
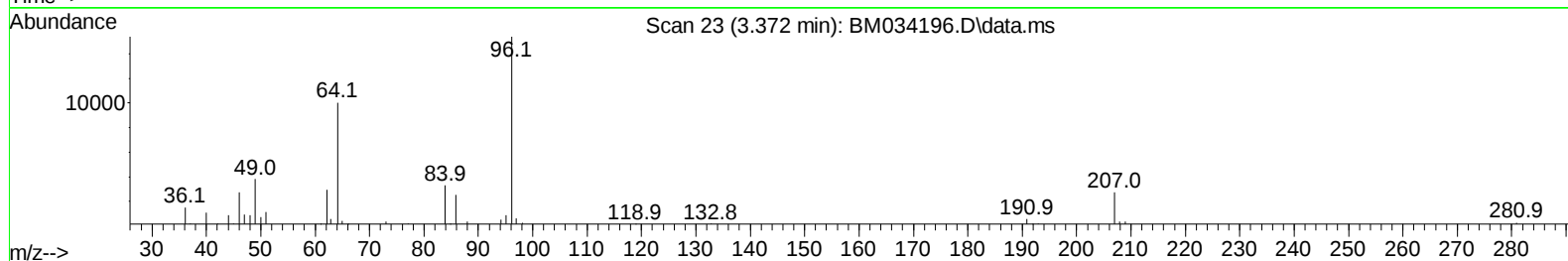
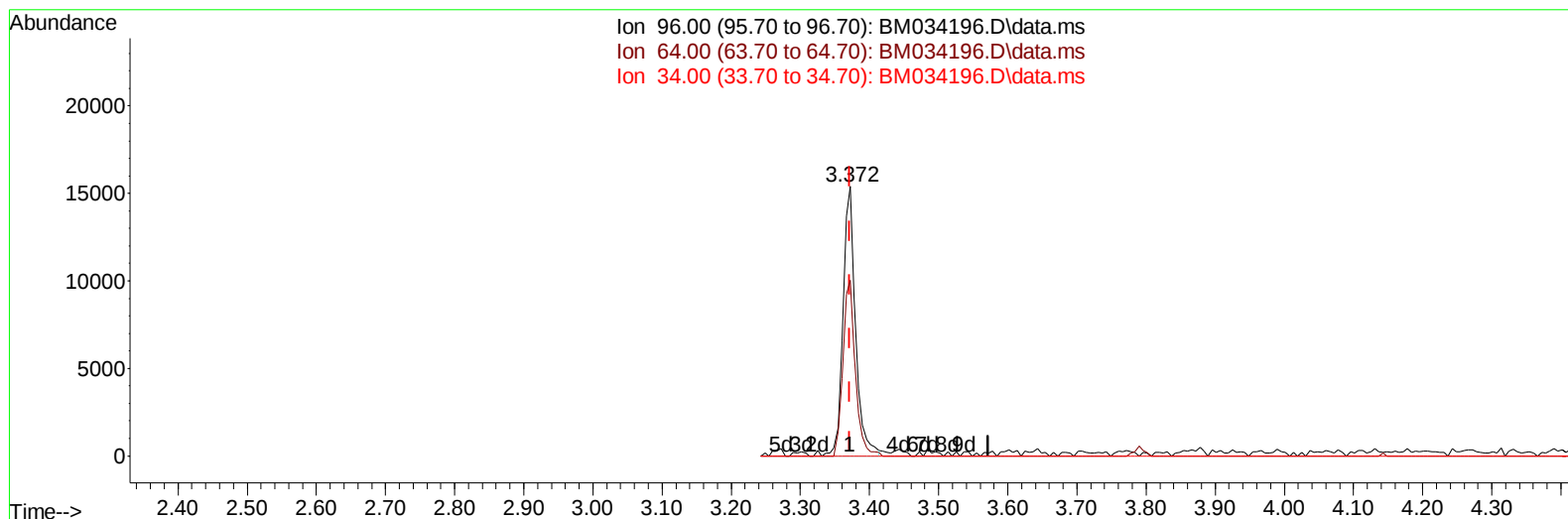
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TIC: BM034196.D\data.ms

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

3.372min (+ 0.000) 5.79 ng/uL m

response 19401

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	65.12
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	154121	20.000	ng/ul	0.00
20) Naphthalene-d8	10.801	136	615129	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.624	164	408975	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.365	188	765594	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.530	240	631418	20.000	ng/ul	-0.01
88) Perylene-d12	23.977	264	683892	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	19401m	5.790	ng/uL	0.00
4) Pyridine-d5	3.790	84	251706	25.804	ng/ul	0.00
7) Phenol-d5	7.137	99	322701	26.722	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.307	67	185633	28.559	ng/ul	0.00
11) 2-Chlorophenol-d4	7.513	132	285209	28.491	ng/ul	0.00
15) 4-Methylphenol-d8	8.689	113	272742	26.954	ng/ul	0.00
21) Nitrobenzene-d5	9.148	128	140130	29.401	ng/ul	0.00
24) 2-Nitrophenol-d4	9.878	143	156505	30.234	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.413	165	290894	27.250	ng/ul	0.00
31) 4-Chloroaniline-d4	10.930	131	387695	26.581	ng/ul	0.00
46) Dimethylphthalate-d6	14.030	166	924625	29.794	ng/ul	0.00
49) Acenaphthylene-d8	14.318	160	1114393	28.715	ng/ul	0.00
54) 4-Nitrophenol-d4	14.795	143	125564	21.744	ng/ul	-0.01
60) Fluorene-d10	15.612	176	773494	28.004	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.718	200	118481	23.981	ng/ul	0.00
73) Anthracene-d10	17.459	188	1124680	29.822	ng/ul	0.00
81) Pyrene-d10	19.747	212	1167610	33.642	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.818	264	1121066	30.902	ng/ul	-0.01

Target Compounds Qvalue

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

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