

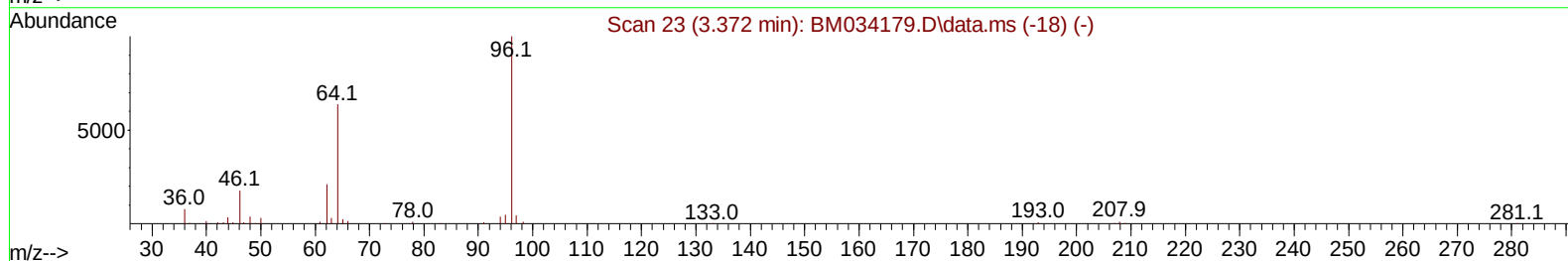
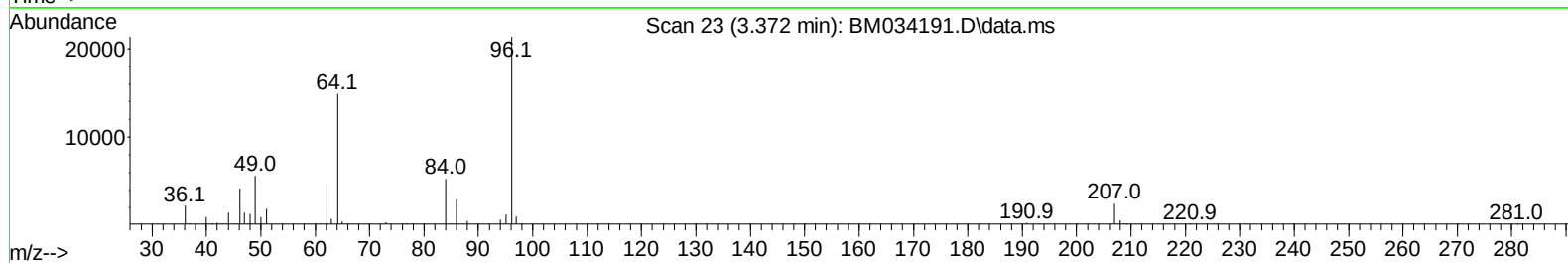
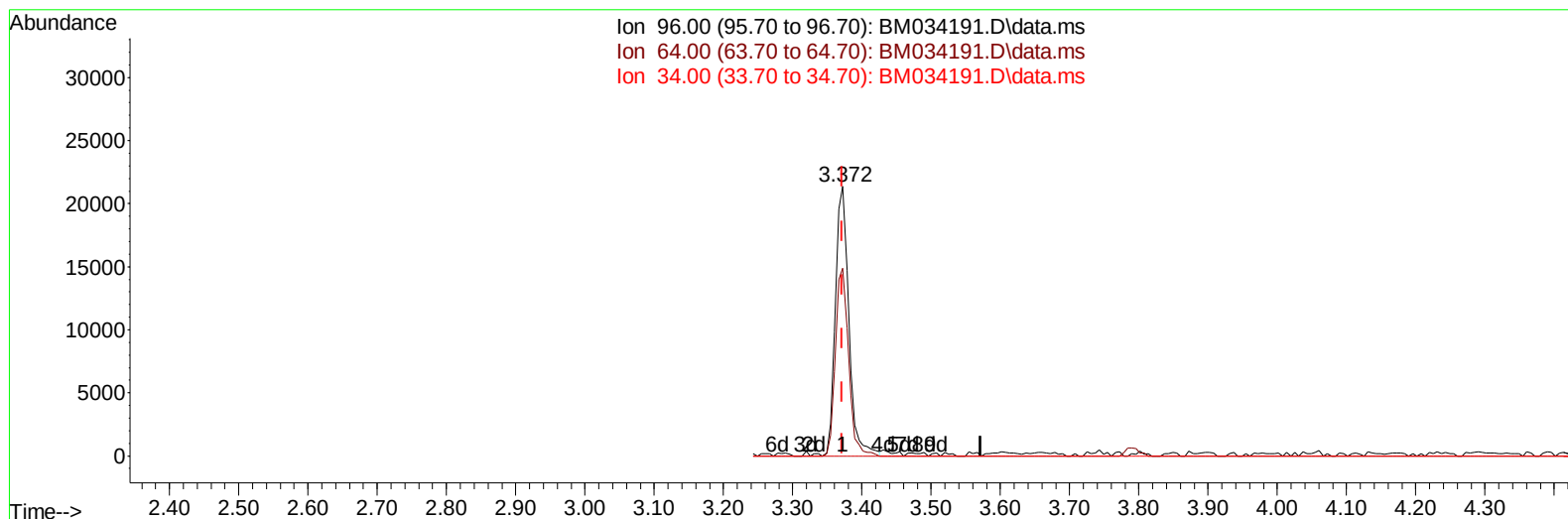
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034191.D
 Acq On : 09 Mar 2022 18:10
 Operator : CG/JU
 Sample : N1775-18
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBSC1

Manual IntegrationsAPPROVED

Quant Time: Mar 10 00:00:28 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: BM034191.D\data.ms

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

3.372min (+ 0.000) 8.45 ng/uL

response 28812

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	69.59
34.00	0.00	0.00
0.00	0.00	0.00

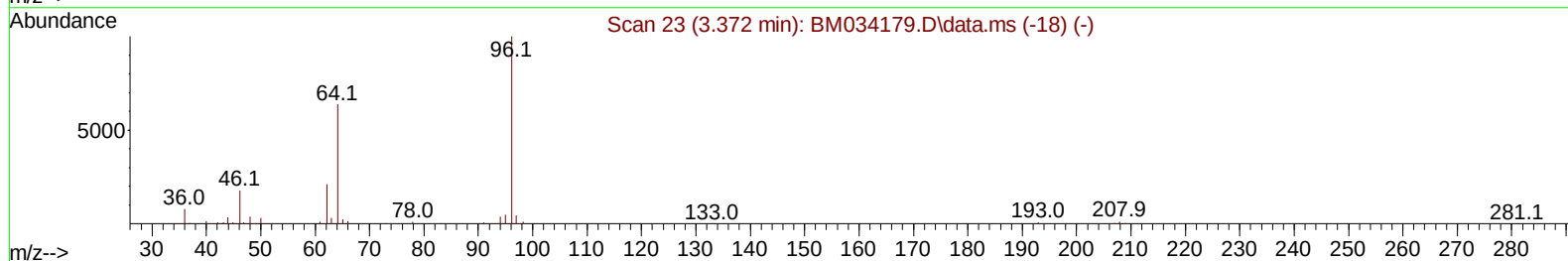
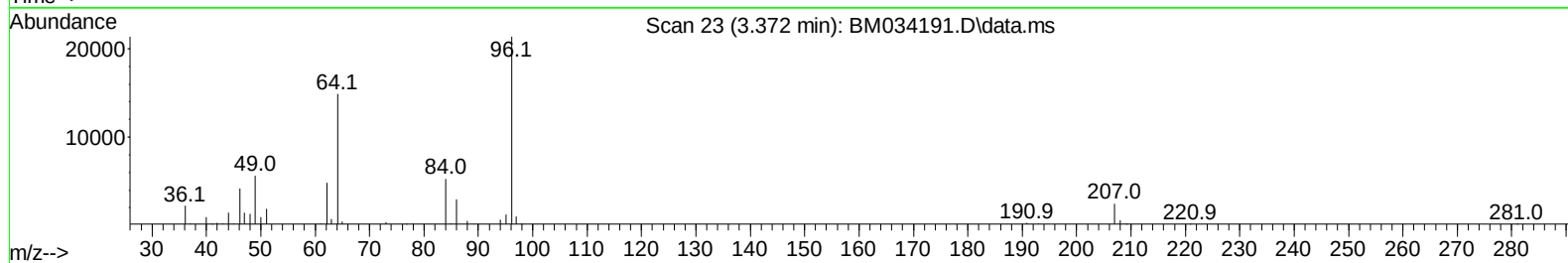
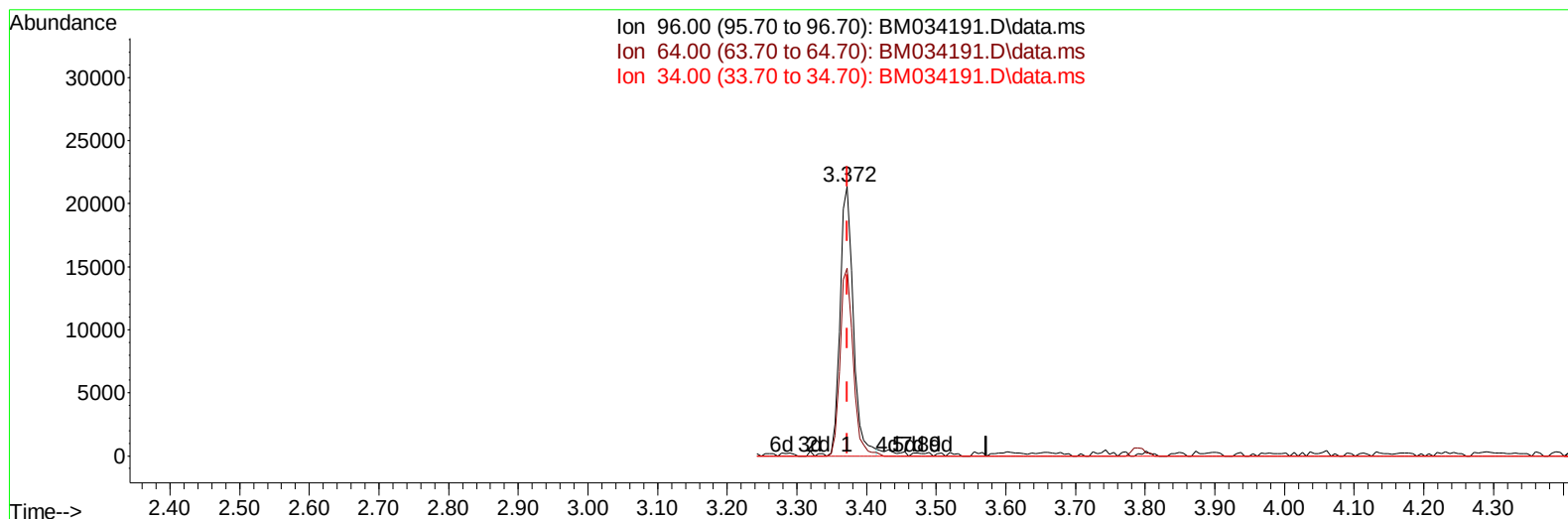
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TIC: BM034191.D\data.ms

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

3.372min (+ 0.000) 8.37 ng/uL m

response 28538

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	69.59
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	7.990	152	156789	20.000	ng/ul	0.00
20) Naphthalene-d8	10.801	136	631790	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.625	164	417583	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.366	188	815653	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.536	240	725874	20.000	ng/ul	# 0.00
88) Perylene-d12	23.983	264	790350	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	28538m	8.371	ng/uL	0.00
4) Pyridine-d5	3.790	84	385121	38.809	ng/ul	0.00
7) Phenol-d5	7.137	99	537563	43.756	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.307	67	310499	46.957	ng/ul	0.00
11) 2-Chlorophenol-d4	7.513	132	473733	46.519	ng/ul	0.00
15) 4-Methylphenol-d8	8.690	113	470758	45.732	ng/ul	0.00
21) Nitrobenzene-d5	9.148	128	240943	49.220	ng/ul	0.00
24) 2-Nitrophenol-d4	9.878	143	276271	51.963	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.419	165	515371	47.004	ng/ul	0.00
31) 4-Chloroaniline-d4	10.931	131	665541	44.428	ng/ul	0.00
46) Dimethylphthalate-d6	14.030	166	1610769	50.833	ng/ul	0.00
49) Acenaphthylene-d8	14.319	160	1984235	50.075	ng/ul	0.00
54) 4-Nitrophenol-d4	14.801	143	232286	39.396	ng/ul	0.00
60) Fluorene-d10	15.619	176	1379311	48.909	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.724	200	227744	43.267	ng/ul	0.00
73) Anthracene-d10	17.466	188	2037941	50.722	ng/ul	0.00
81) Pyrene-d10	19.754	212	2222804	55.710	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.824	264	2161463	51.554	ng/ul	0.00

Target Compounds Qvalue

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

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