

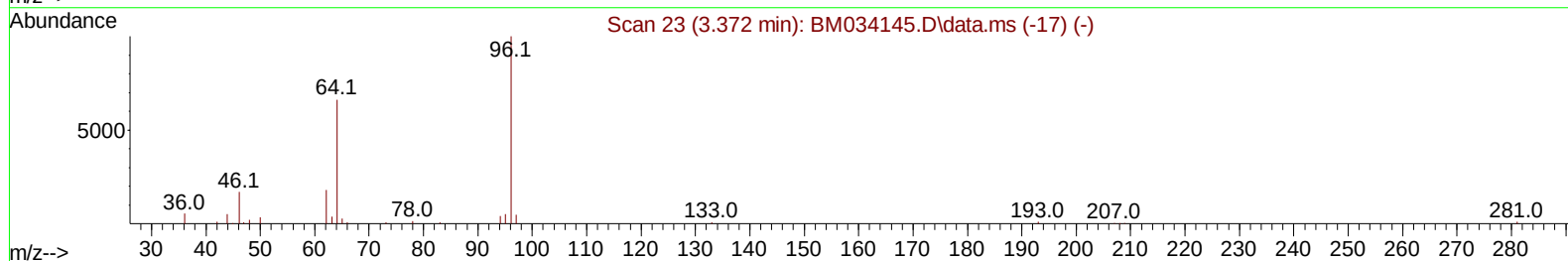
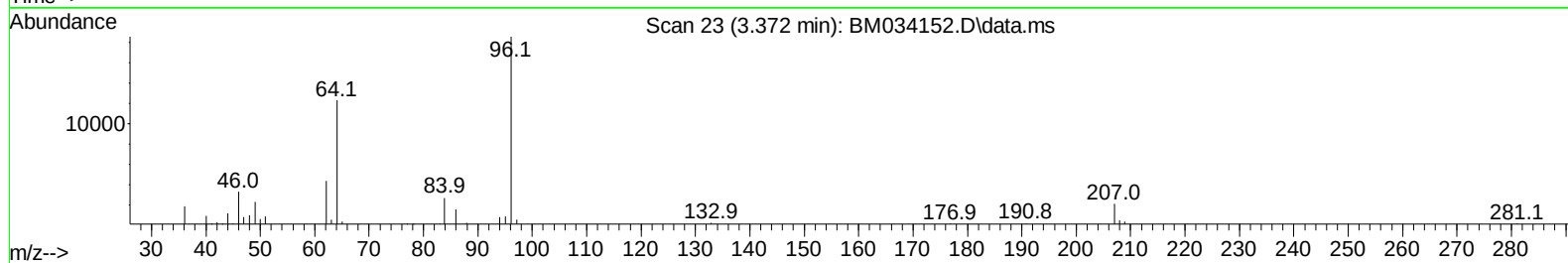
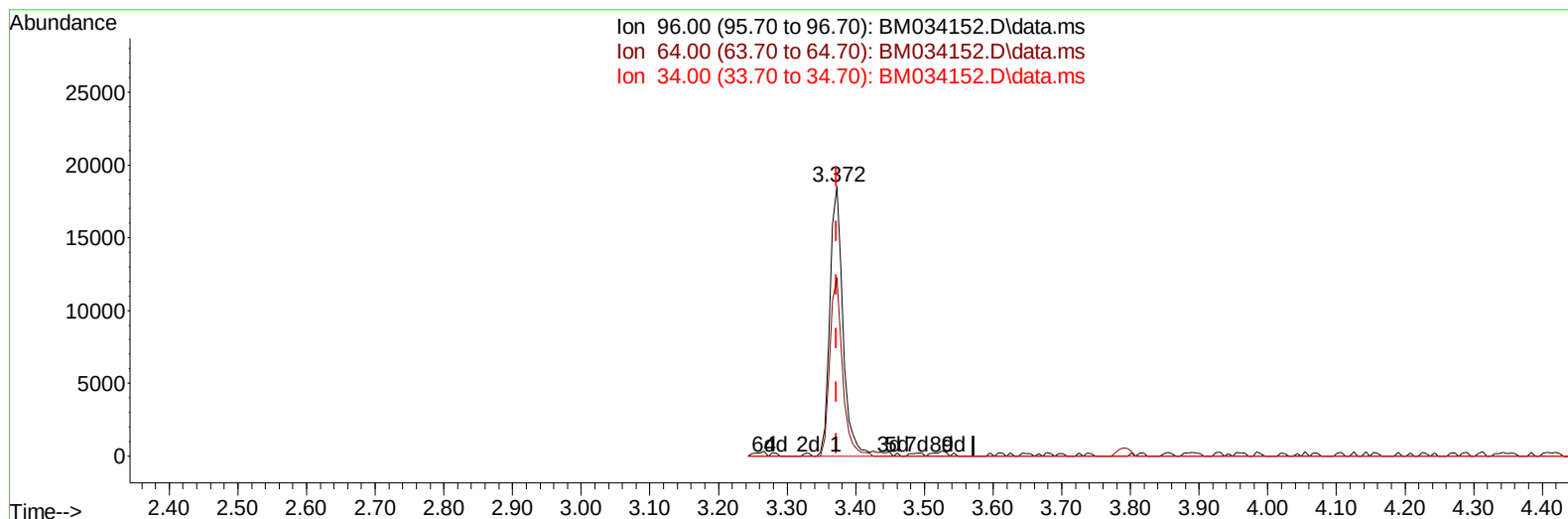
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034152.D
 Acq On : 08 Mar 2022 16:48
 Operator : CG/JU
 Sample : N1733-04
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBRR8

Manual IntegrationsAPPROVED

Quant Time: Mar 09 00:31:07 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 03/09/2022
 Supervised By :Yogesh Patel 03/09/2022



TIC: BM034152.D\data.ms

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

3.372min (+ 0.000) 7.33 ng/uL

response 24909

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	66.25
34.00	0.00	0.00
0.00	0.00	0.00

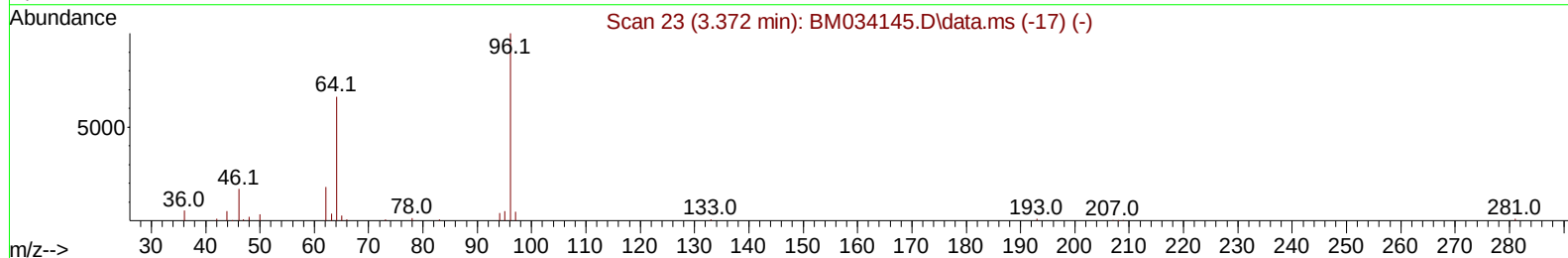
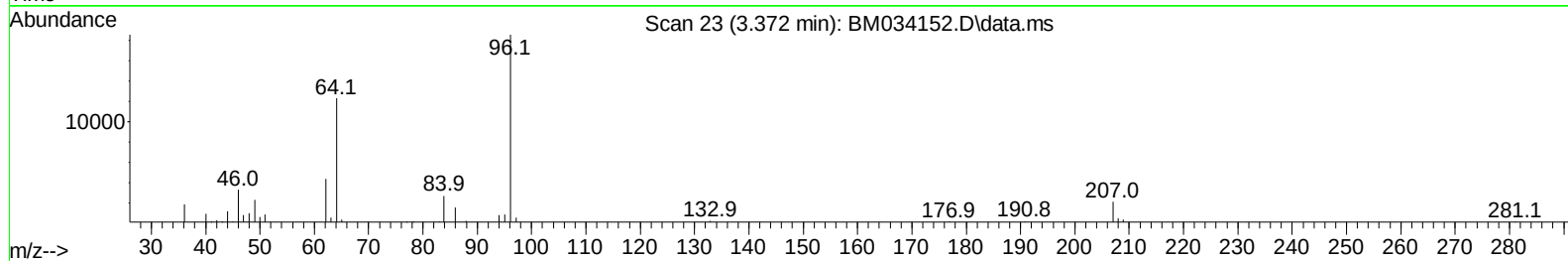
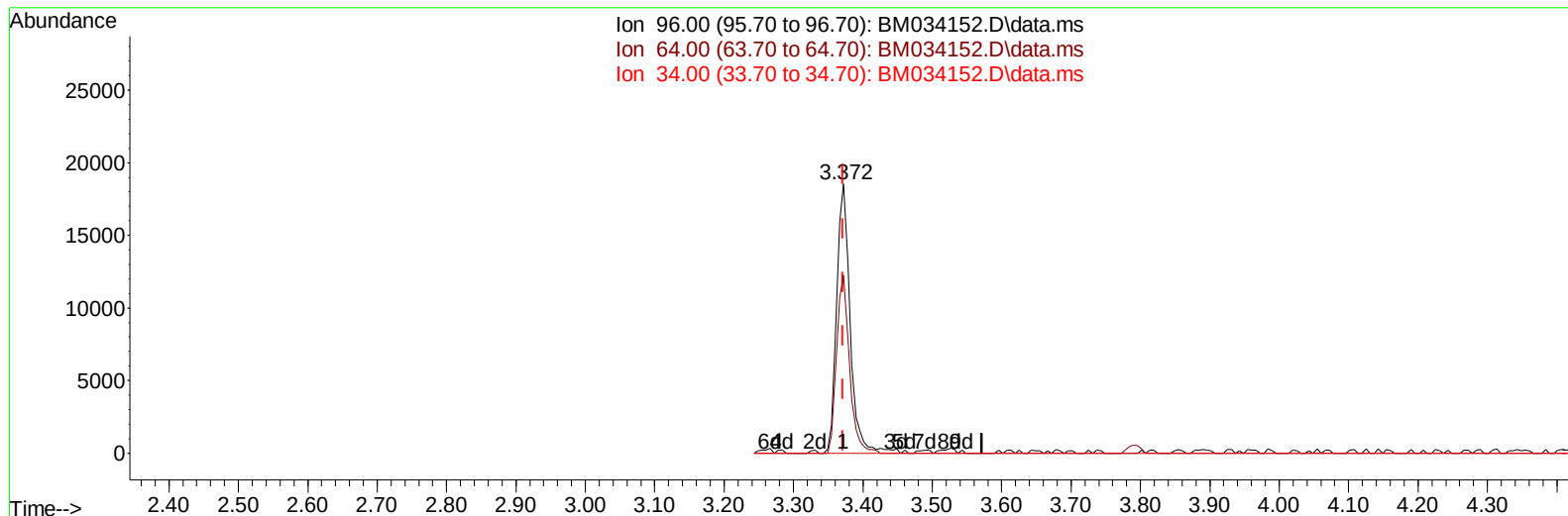
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(3) 1,4-Dioxane-d8 (S)

3.372min (+ 0.000) 7.22 ng/uL m

response 24553

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	66.25
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	7.995	152	156361	20.000	ng/ul	0.00
20) Naphthalene-d8	10.807	136	631029	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.630	164	430229	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.371	188	906271	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.542	240	839107	20.000	ng/ul	0.00
88) Perylene-d12	23.994	264	802507	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	24553m	7.222	ng/uL	0.00
4) Pyridine-d5	3.790	84	316284	31.960	ng/ul	0.00
7) Phenol-d5	7.143	99	435707	35.563	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.313	67	247068	37.467	ng/ul	0.00
11) 2-Chlorophenol-d4	7.519	132	381377	37.552	ng/ul	0.00
15) 4-Methylphenol-d8	8.695	113	373037	36.338	ng/ul	0.00
21) Nitrobenzene-d5	9.154	128	193473	39.571	ng/ul	0.00
24) 2-Nitrophenol-d4	9.884	143	214969	40.481	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.425	165	397342	36.283	ng/ul	0.00
31) 4-Chloroaniline-d4	10.936	131	530248	35.439	ng/ul	0.00
46) Dimethylphthalate-d6	14.036	166	1306601	40.022	ng/ul	0.00
49) Acenaphthylene-d8	14.325	160	1594994	39.069	ng/ul	0.00
54) 4-Nitrophenol-d4	14.807	143	179721	29.585	ng/ul	0.00
60) Fluorene-d10	15.619	176	1117996	38.477	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.724	200	176188	30.125	ng/ul	0.00
73) Anthracene-d10	17.471	188	1749757	39.195	ng/ul	0.00
81) Pyrene-d10	19.759	212	1988537	43.113	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.836	264	1689590	39.689	ng/ul	0.00

Target Compounds Qvalue

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

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