

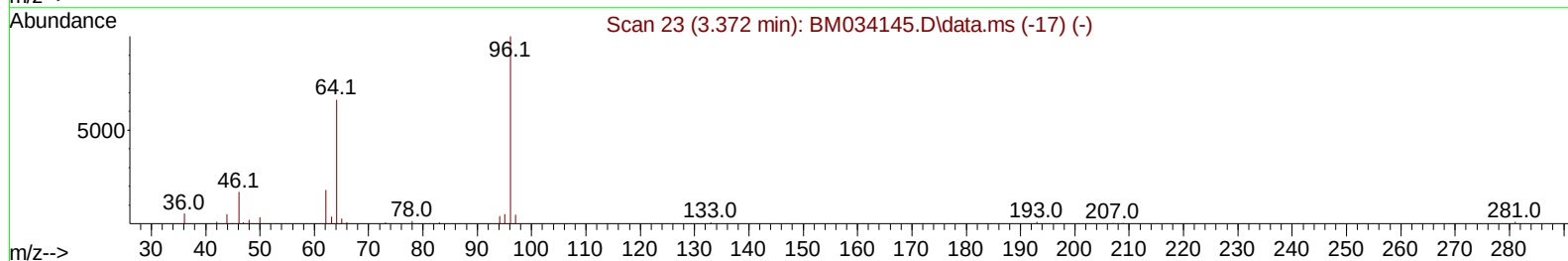
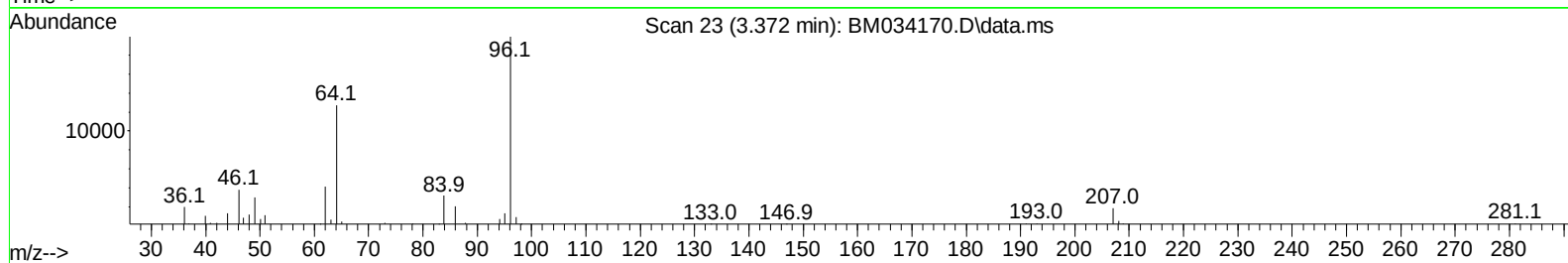
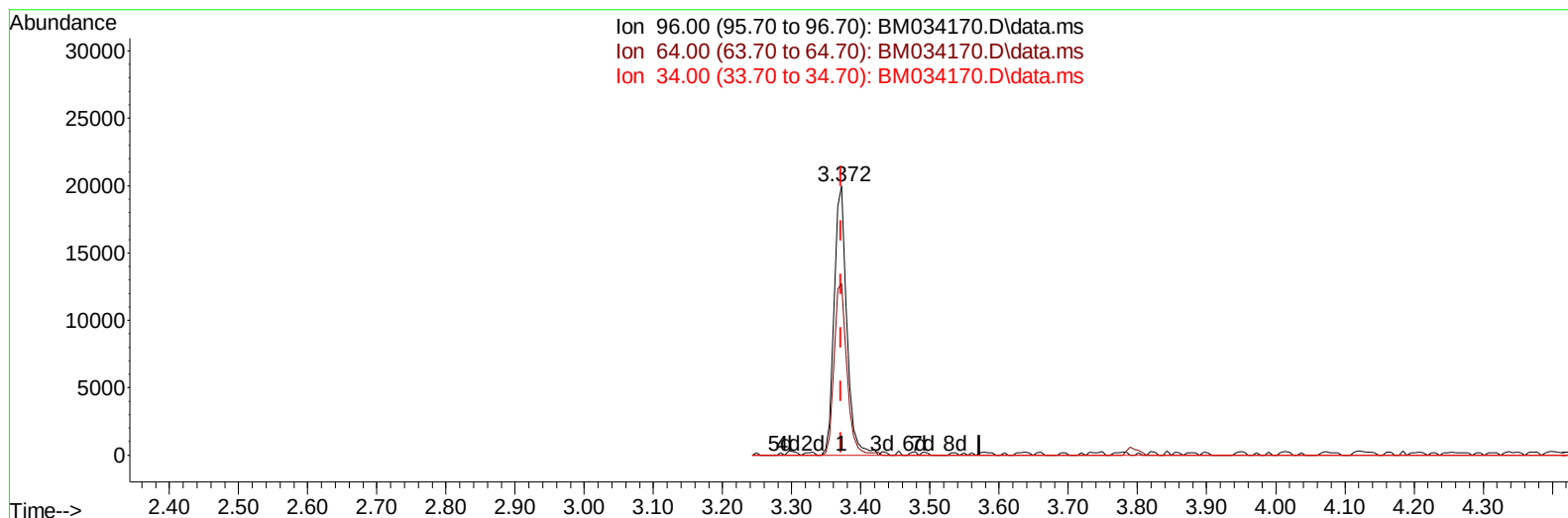
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034170.D
 Acq On : 09 Mar 2022 04:19
 Operator : CG/JU
 Sample : N1733-07
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBRS2

Manual IntegrationsAPPROVED

Quant Time: Mar 09 05:10:53 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: BM034170.D\data.ms

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

3.372min (+ 0.000) 8.08 ng/uL

response 25245

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	63.78
34.00	0.00	0.00
0.00	0.00	0.00

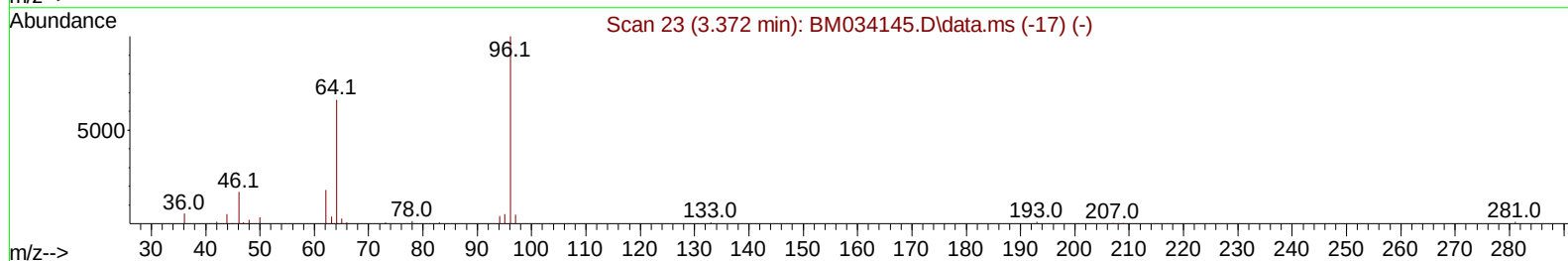
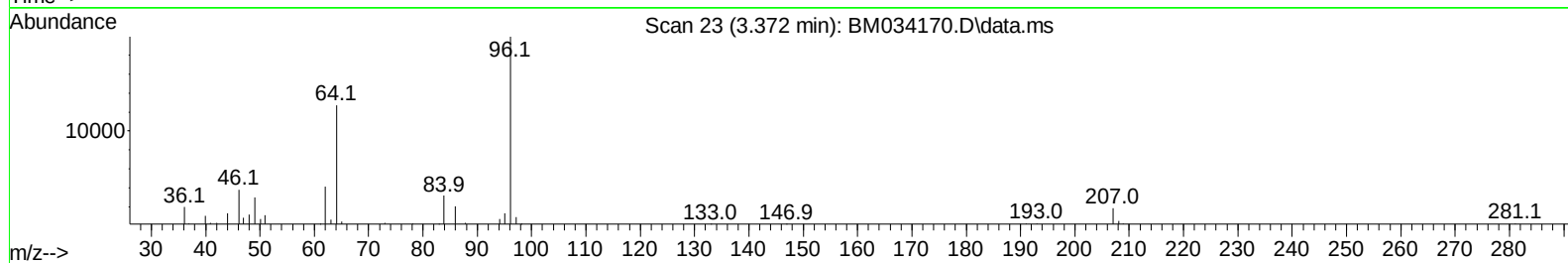
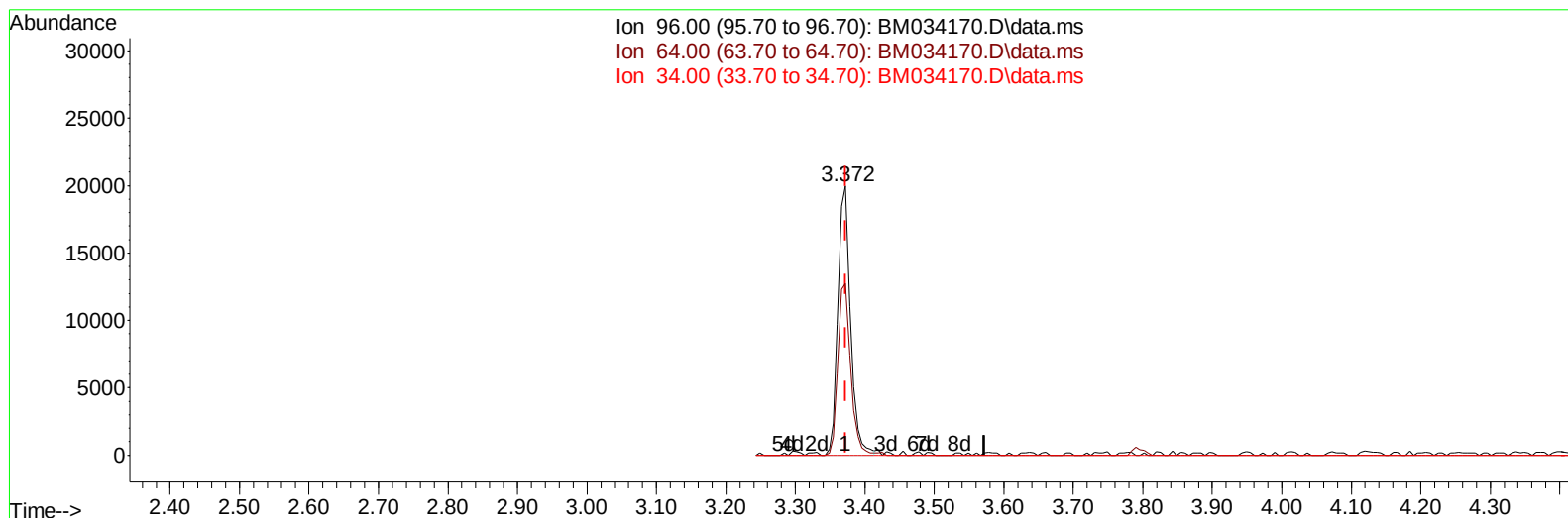
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TIC: BM034170.D\data.ms

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

3.372min (+ 0.000) 8.03 ng/uL m

response 25108

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	63.78
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	143724	20.000	ng/ul	0.00
20) Naphthalene-d8	10.801	136	553447	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.624	164	340941	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.365	188	676062	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.536	240	674986	20.000	ng/ul	# 0.00
88) Perylene-d12	23.977	264	754601	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	25108m	8.035	ng/uL	0.00
4) Pyridine-d5	3.790	84	285816	31.420	ng/ul	0.00
7) Phenol-d5	7.137	99	428030	38.008	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.307	67	244703	40.371	ng/ul	0.00
11) 2-Chlorophenol-d4	7.519	132	389283	41.701	ng/ul	0.00
15) 4-Methylphenol-d8	8.690	113	360764	38.232	ng/ul	0.00
21) Nitrobenzene-d5	9.148	128	185614	43.285	ng/ul	0.00
24) 2-Nitrophenol-d4	9.878	143	209478	44.977	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.419	165	386536	40.244	ng/ul	0.00
31) 4-Chloroaniline-d4	10.931	131	483169	36.819	ng/ul	0.00
46) Dimethylphthalate-d6	14.030	166	1094678	42.312	ng/ul	0.00
49) Acenaphthylene-d8	14.319	160	1379613	42.643	ng/ul	0.00
54) 4-Nitrophenol-d4	14.801	143	163278	33.918	ng/ul	0.00
60) Fluorene-d10	15.618	176	948905	41.211	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.724	200	153945	35.285	ng/ul	0.00
73) Anthracene-d10	17.465	188	1431372	42.981	ng/ul	0.00
81) Pyrene-d10	19.754	212	1614661	43.519	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.824	264	1755810	43.863	ng/ul	0.00

Target Compounds Qvalue

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

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