

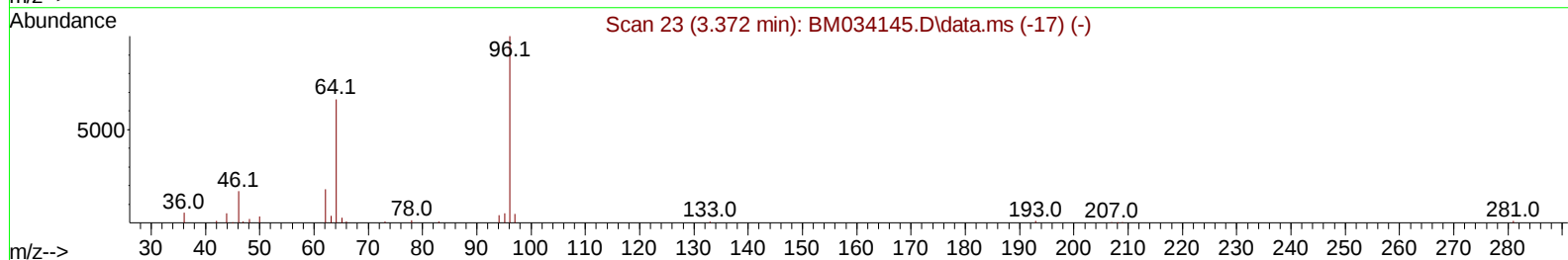
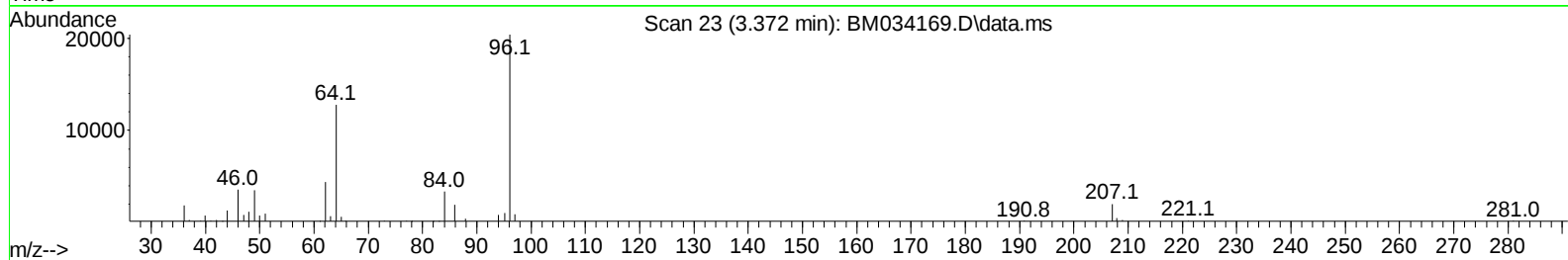
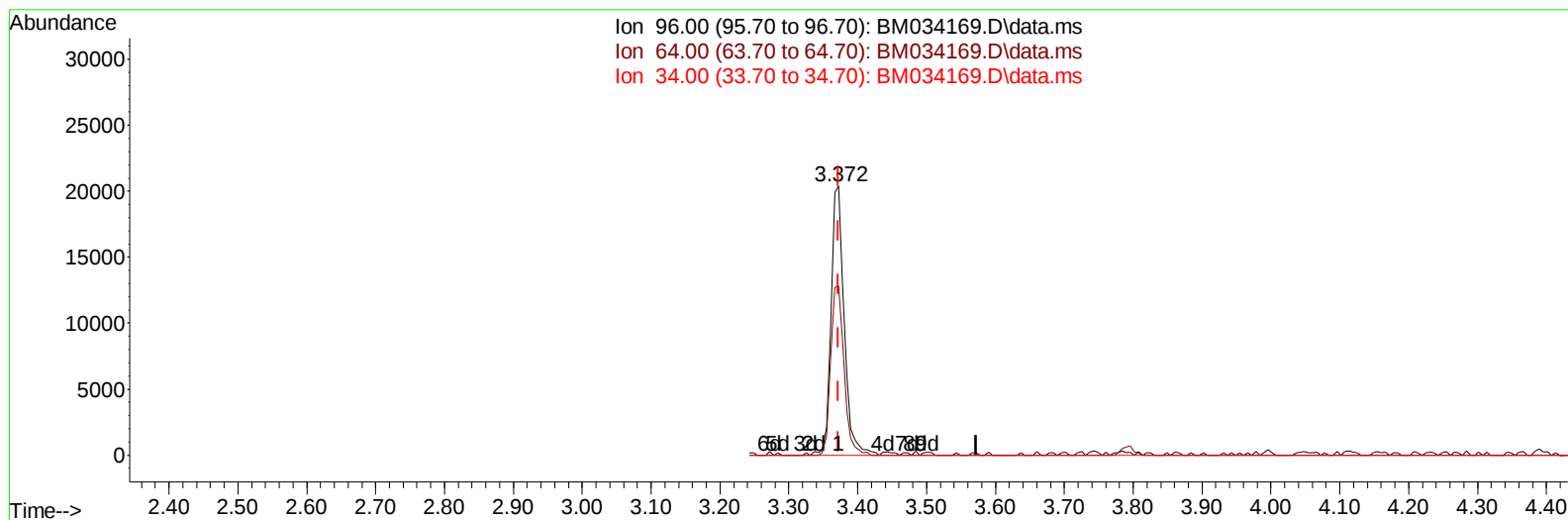
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034169.D
 Acq On : 09 Mar 2022 03:43
 Operator : CG/JU
 Sample : N1733-14
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DBRS7

Manual IntegrationsAPPROVED

Quant Time: Mar 09 04:22:36 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: BM034169.D\data.ms

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

3.372min (+ 0.000) 8.01 ng/uL

response 26903

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	62.83
34.00	0.00	0.00
0.00	0.00	0.00

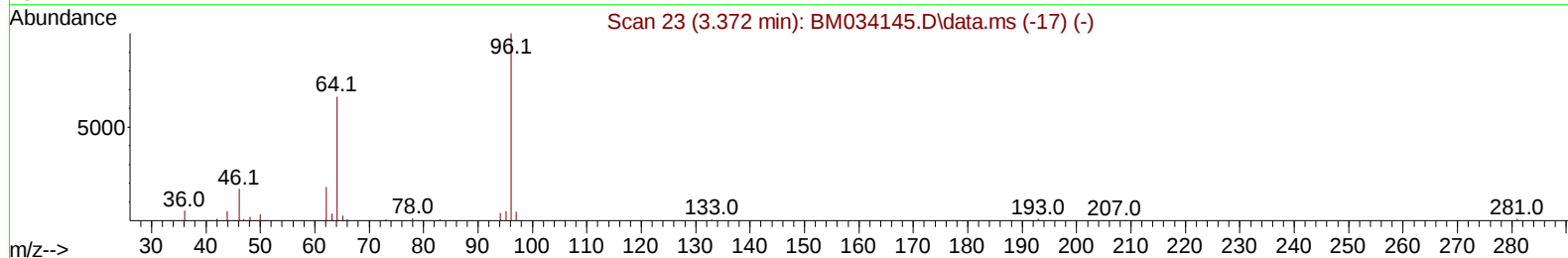
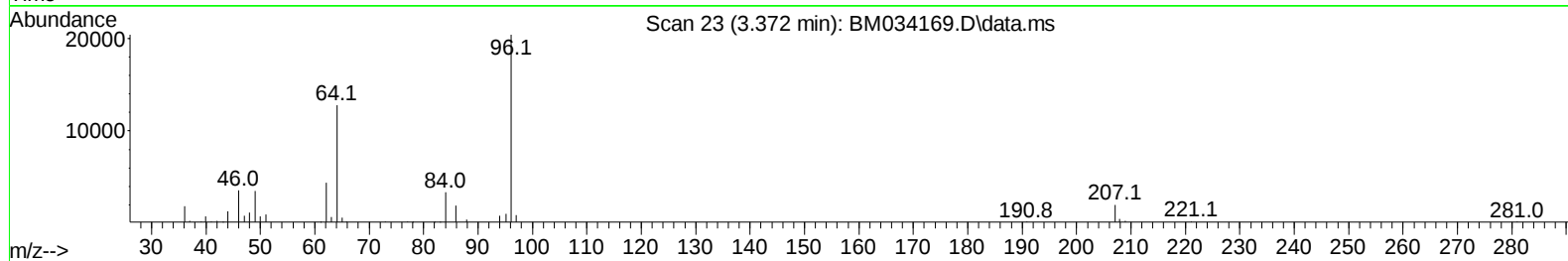
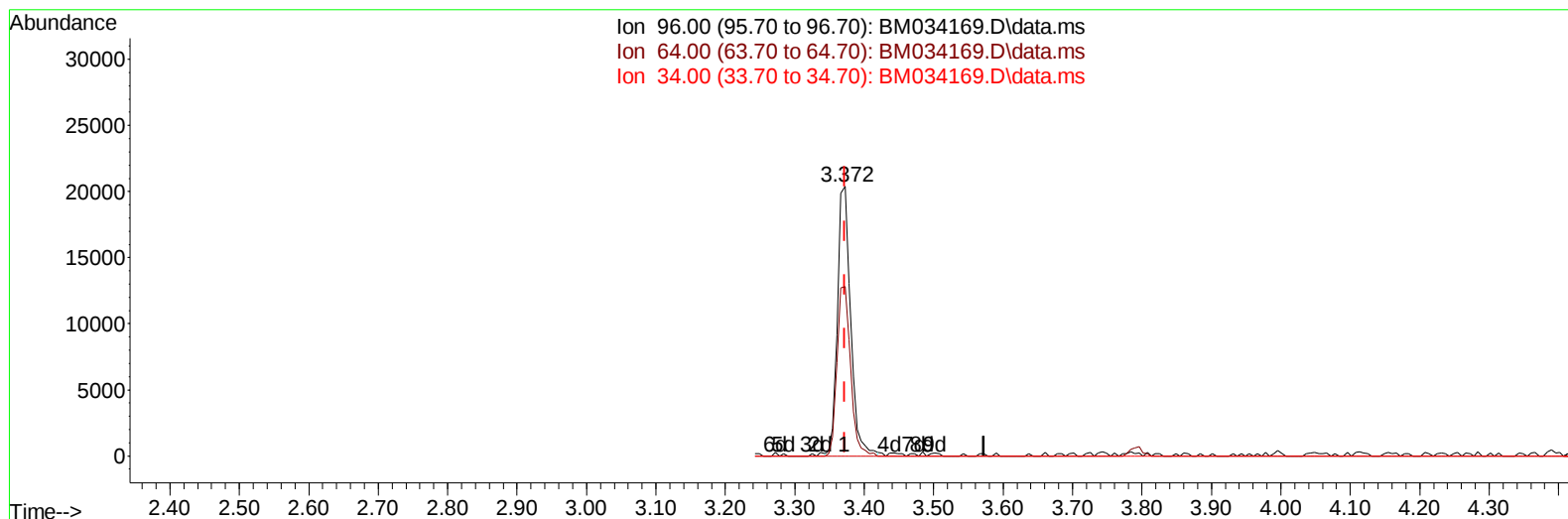
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TIC: BM034169.D\data.ms

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

3.372min (+ 0.000) 7.98 ng/uL m

response 26822

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	62.83
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	154524	20.000	ng/ul	0.00
20) Naphthalene-d8	10.801	136	619796	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.624	164	423351	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.365	188	847154	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.536	240	701929	20.000	ng/ul	# 0.00
88) Perylene-d12	23.983	264	726578	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	26822m	7.983	ng/uL	0.00
4) Pyridine-d5	3.790	84	324201	33.149	ng/ul	0.00
7) Phenol-d5	7.143	99	499460	41.251	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.313	67	282971	43.421	ng/ul	0.00
11) 2-Chlorophenol-d4	7.519	132	435741	43.415	ng/ul	0.00
15) 4-Methylphenol-d8	8.690	113	433312	42.711	ng/ul	0.00
21) Nitrobenzene-d5	9.154	128	223128	46.463	ng/ul	0.00
24) 2-Nitrophenol-d4	9.878	143	255991	49.080	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.419	165	474942	44.155	ng/ul	0.00
31) 4-Chloroaniline-d4	10.931	131	620263	42.206	ng/ul	0.00
46) Dimethylphthalate-d6	14.030	166	1530939	47.655	ng/ul	0.00
49) Acenaphthylene-d8	14.319	160	1888256	47.004	ng/ul	0.00
54) 4-Nitrophenol-d4	14.801	143	206088	34.477	ng/ul	0.00
60) Fluorene-d10	15.619	176	1330333	46.529	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.724	200	196725	35.984	ng/ul	0.00
73) Anthracene-d10	17.465	188	1984749	47.561	ng/ul	0.00
81) Pyrene-d10	19.754	212	2067460	53.585	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.824	264	1868104	48.468	ng/ul	0.00

Target Compounds Qvalue

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

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