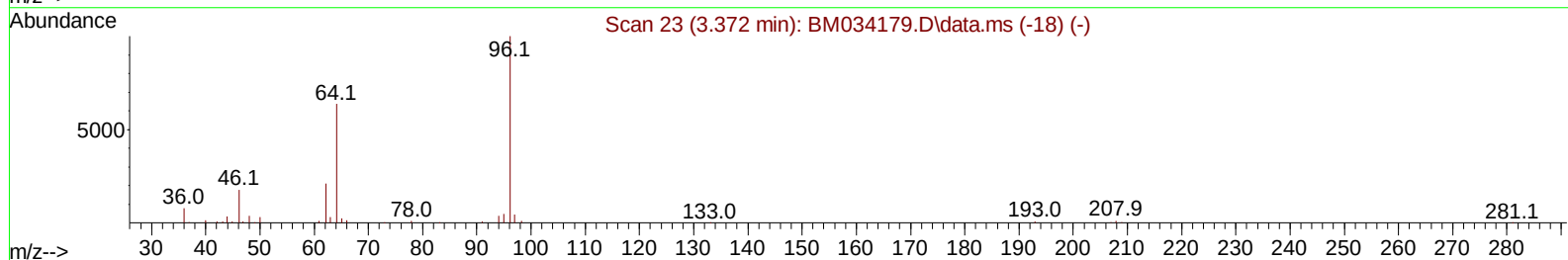
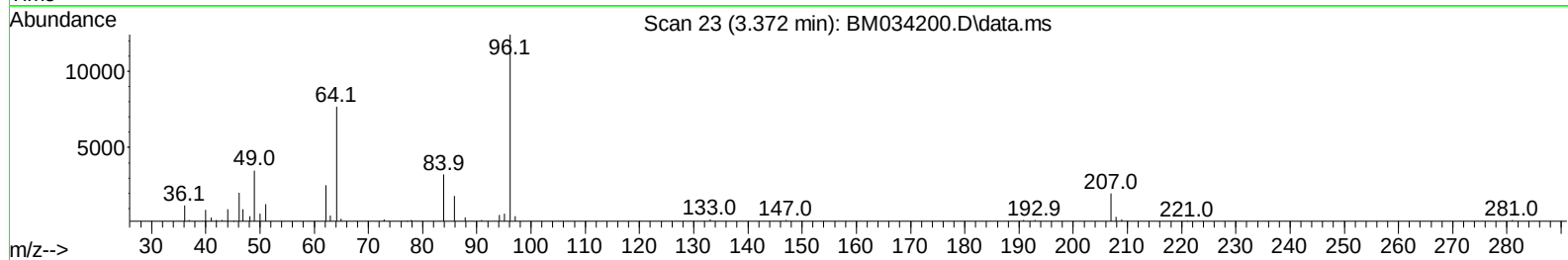
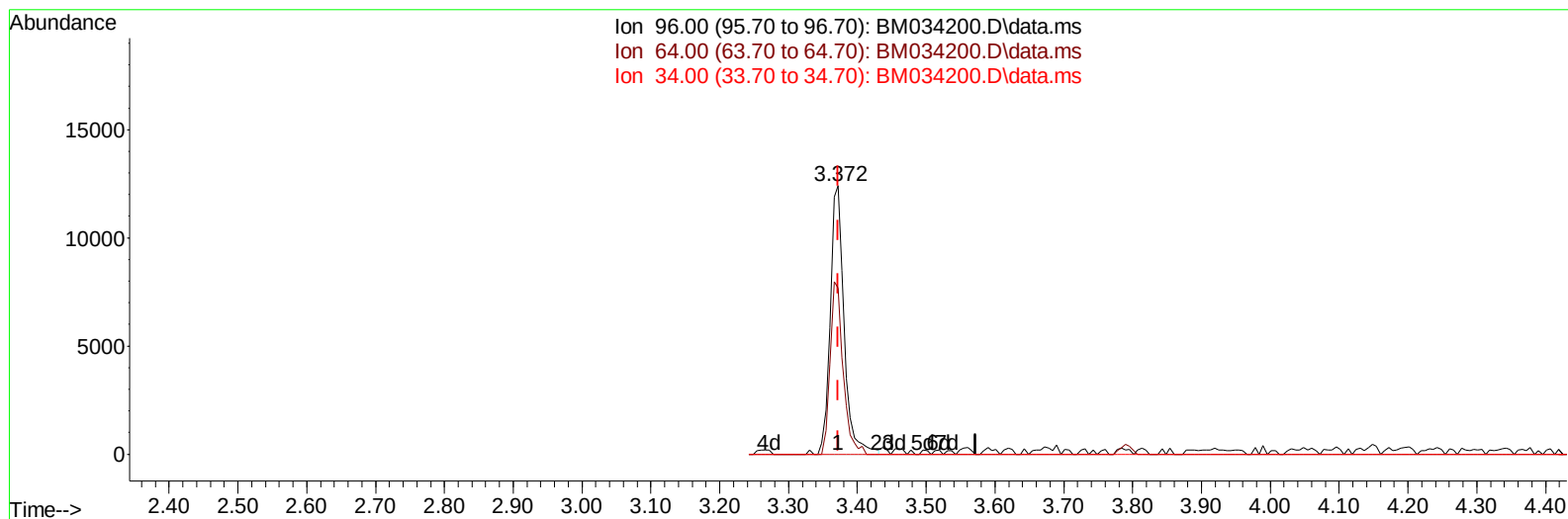


Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
Data File : BM034200.D
Acq On : 10 Mar 2022 00:12
Operator : CG/JU
Sample : N1775-27
Misc :
ALS Vial : 60 Sample Multiplier: 1

Quant Time: Mar 10 01:04:34 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



TIC: BM034200.D\data.ms

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

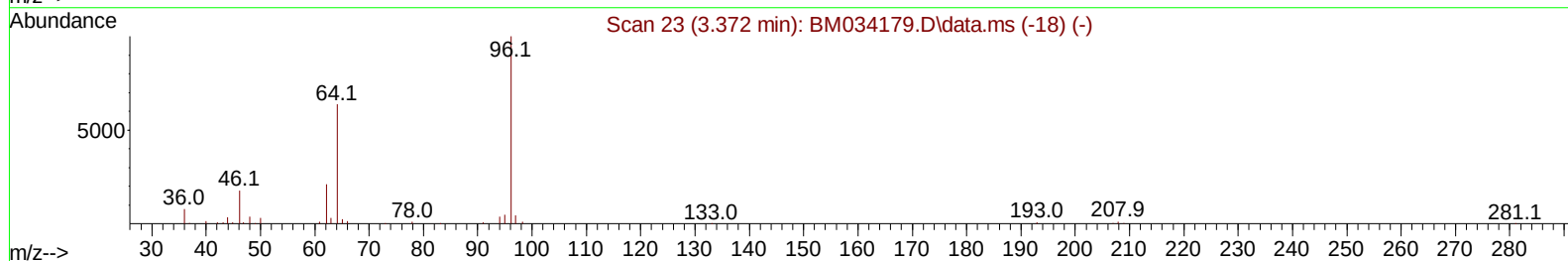
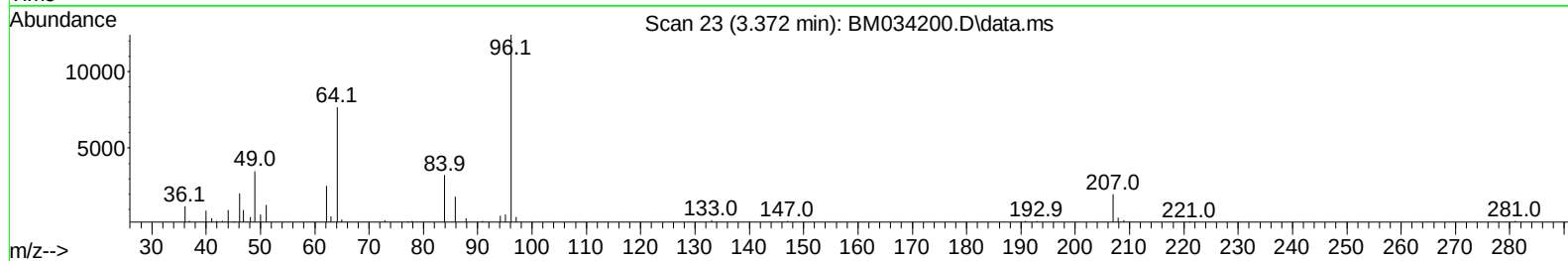
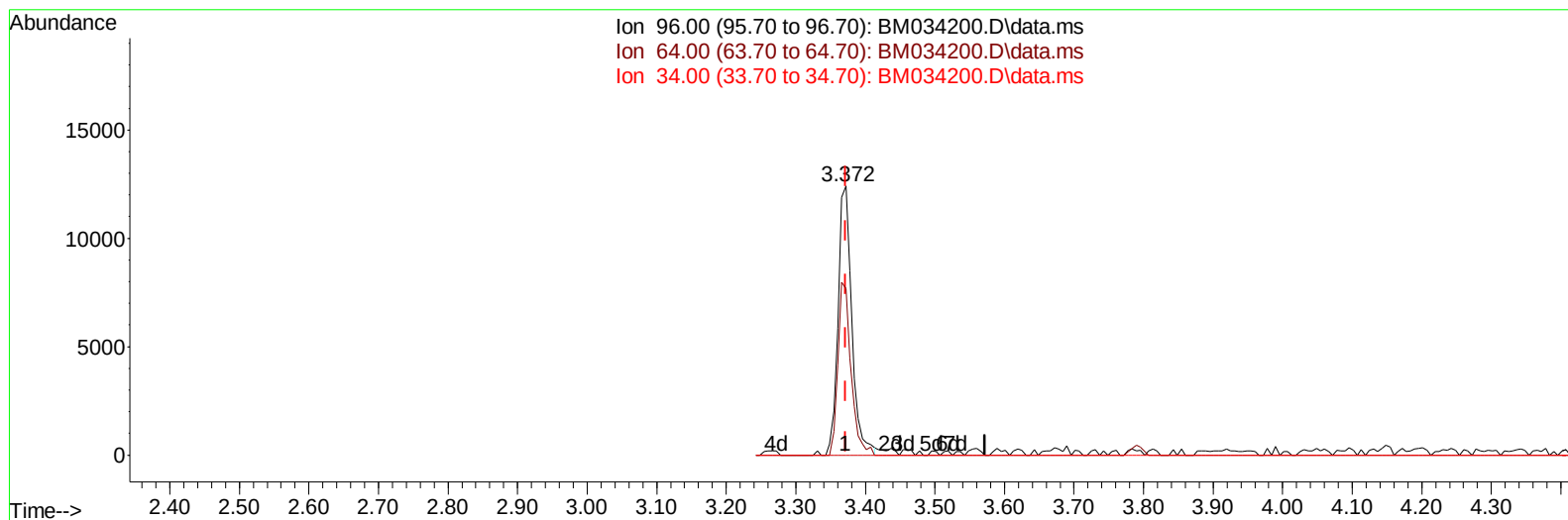
3.372min (0.000) 5.74 ng/uL

response 17362

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	61.97
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
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Operator : CG/JU
Sample : N1775-27
Misc :
ALS Vial : 60 Sample Multiplier: 1

Quant Time: Mar 10 01:04:34 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM030822.M
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TIC: BM034200.D\data.ms

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

3.372min (0.000) 5.66 ng/uL m

response 17120

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.20	61.97
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM030822\
 Data File : BM034200.D
 Acq On : 10 Mar 2022 00:12
 Operator : CG/JU
 Sample : N1775-27
 Misc :
 ALS Vial : 60 Sample Multiplier: 1

Quant Time: Mar 10 01:04:34 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

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.983	152	139016	20.000	ng/ul	-0.01
20) Naphthalene-d8	10.801	136	560656	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.624	164	370193	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.365	188	708981	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.530	240	598409	20.000	ng/ul	#-0.01
88) Perylene-d12	23.971	264	644256	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	17120m	5.664	ng/uL	0.00
4) Pyridine-d5	3.790	84	205299	23.333	ng/ul	0.00
7) Phenol-d5	7.136	99	297400	27.303	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.307	67	172877	29.487	ng/ul	0.00
11) 2-Chlorophenol-d4	7.513	132	260499	28.850	ng/ul	0.00
15) 4-Methylphenol-d8	8.689	113	255840	28.031	ng/ul	0.00
21) Nitrobenzene-d5	9.148	128	128947	29.684	ng/ul	0.00
24) 2-Nitrophenol-d4	9.877	143	143119	30.334	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.413	165	276642	28.432	ng/ul	0.00
31) 4-Chloroaniline-d4	10.930	131	360375	27.109	ng/ul	0.00
46) Dimethylphthalate-d6	14.030	166	879802	31.319	ng/ul	0.00
49) Acenaphthylene-d8	14.318	160	1076636	30.649	ng/ul	0.00
54) 4-Nitrophenol-d4	14.795	143	87502	16.740	ng/ul	-0.01
60) Fluorene-d10	15.612	176	750248	30.008	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.718	200	84169	18.396	ng/ul	0.00
73) Anthracene-d10	17.459	188	1083321	31.019	ng/ul	0.00
81) Pyrene-d10	19.747	212	1122436	34.124	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.818	264	1065321	31.172	ng/ul	-0.01

Target Compounds Qvalue

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

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