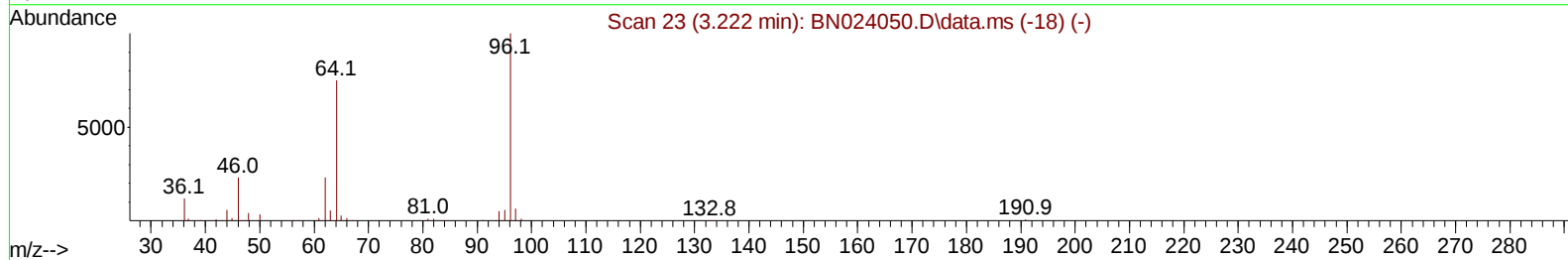
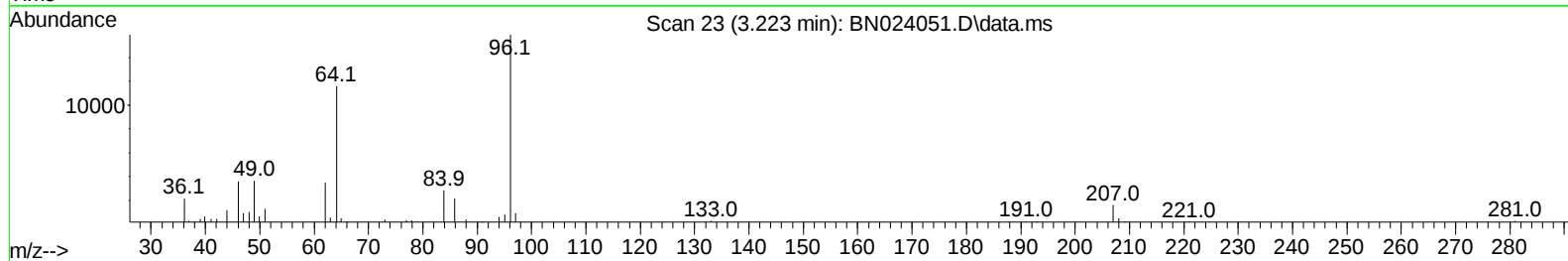
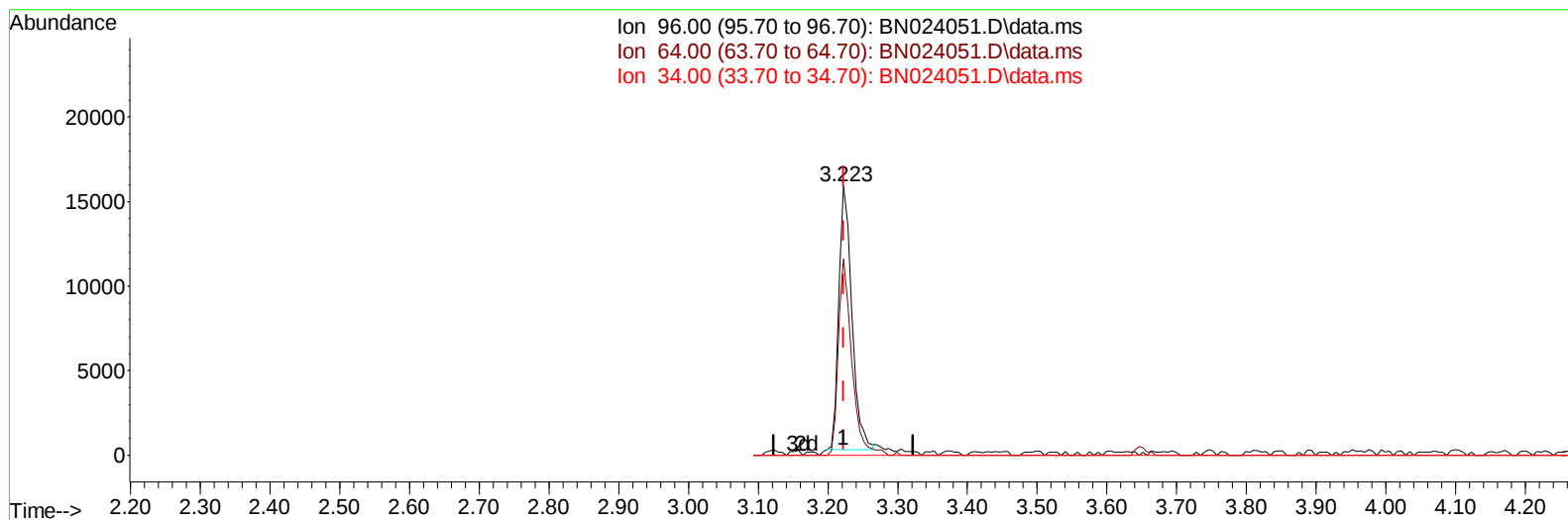


Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021623\
Data File : BN024051.D
Acq On : 16 Feb 2023 13:08
Operator : CG/JU
Sample : PB150871BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Feb 17 05:36:14 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020623.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Feb 17 05:33:41 2023
Response via : Initial Calibration



TIC: BN024051.D\data.ms

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

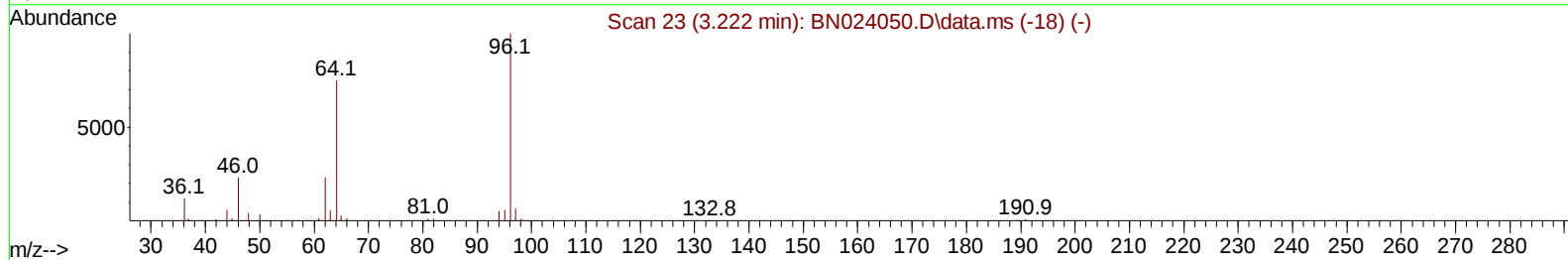
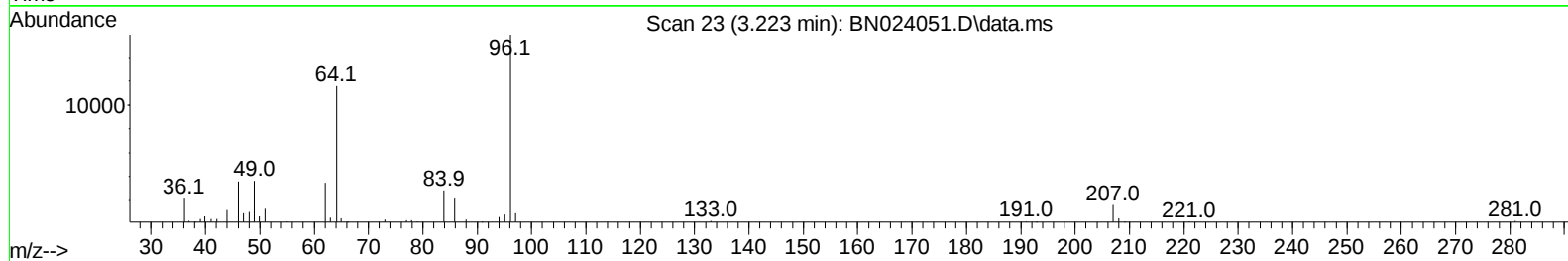
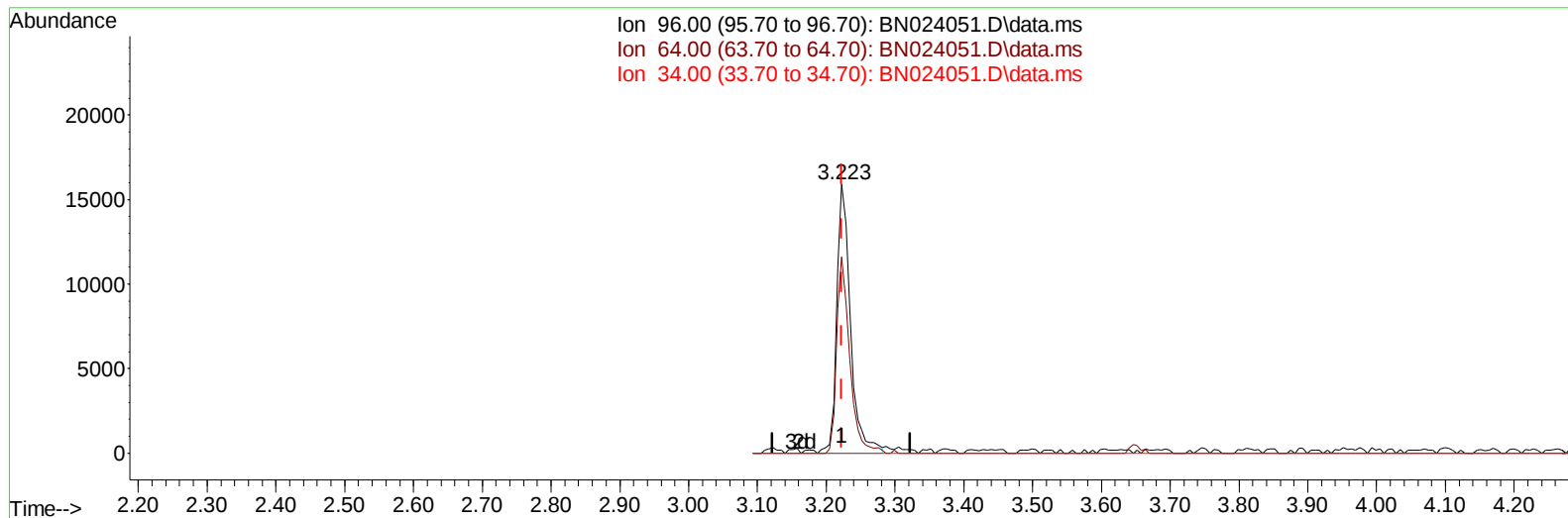
3.223min (+ 0.000) 5.33 ng/uL

response 20271

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.80	72.94
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021623\
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Acq On : 16 Feb 2023 13:08
Operator : CG/JU
Sample : PB150871BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Quant Time: Feb 17 05:36:14 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN020623.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Feb 17 05:33:41 2023
Response via : Initial Calibration



TIC: BN024051.D\data.ms

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

3.223min (+ 0.000) 5.82 ng/uL m

response 22163

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	74.80	72.94
34.00	0.00	0.00
0.00	0.00	0.00

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN021623\
 Data File : BN024051.D
 Acq On : 16 Feb 2023 13:08
 Operator : CG/JU
 Sample : PB150871BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Feb 17 05:36:14 2023
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 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Feb 17 05:33:41 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.828	152	153873	20.000	ng/ul	0.00
20) Naphthalene-d8	10.634	136	667186	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.481	164	442674	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.234	188	1016067	20.000	ng/ul	0.00
79) Chrysene-d12	21.422	240	953481	20.000	ng/ul	0.00
88) Perylene-d12	23.816	264	899863	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.223	96	22163m	5.824	ng/uL	0.00
4) Pyridine-d5	3.646	84	271702	25.740	ng/ul	0.00
7) Phenol-d5	6.999	99	367734	27.956	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.163	67	247545	29.822	ng/ul	0.00
11) 2-Chlorophenol-d4	7.358	132	296925	29.023	ng/ul	0.00
15) 4-Methylphenol-d8	8.546	113	295403	28.746	ng/ul	0.00
21) Nitrobenzene-d5	8.999	128	147210	27.638	ng/ul	0.00
24) 2-Nitrophenol-d4	9.722	143	163900	27.405	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.257	165	292202	26.218	ng/ul	0.00
31) 4-Chloroaniline-d4	10.781	131	415052	27.555	ng/ul	0.00
46) Dimethylphthalate-d6	13.898	166	990076	28.938	ng/ul	0.00
49) Acenaphthylene-d8	14.175	160	1039899	27.435	ng/ul	0.00
54) 4-Nitrophenol-d4	14.692	143	151291	25.706	ng/ul	0.00
60) Fluorene-d10	15.475	176	838968	29.123	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.604	200	149604	21.311	ng/ul	0.00
73) Anthracene-d10	17.334	188	1364688	29.484	ng/ul	0.00
81) Pyrene-d10	19.628	212	1617230	28.157	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.657	264	1387507	30.139	ng/ul	0.01

Target Compounds	Qvalue

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

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