

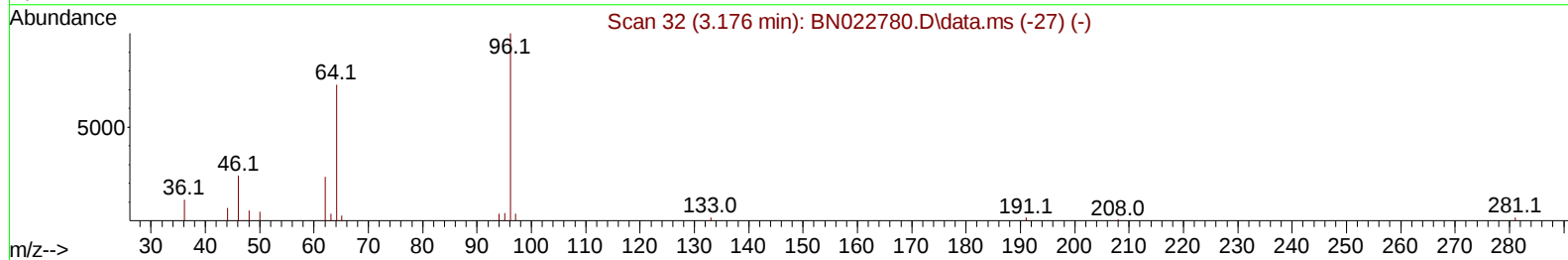
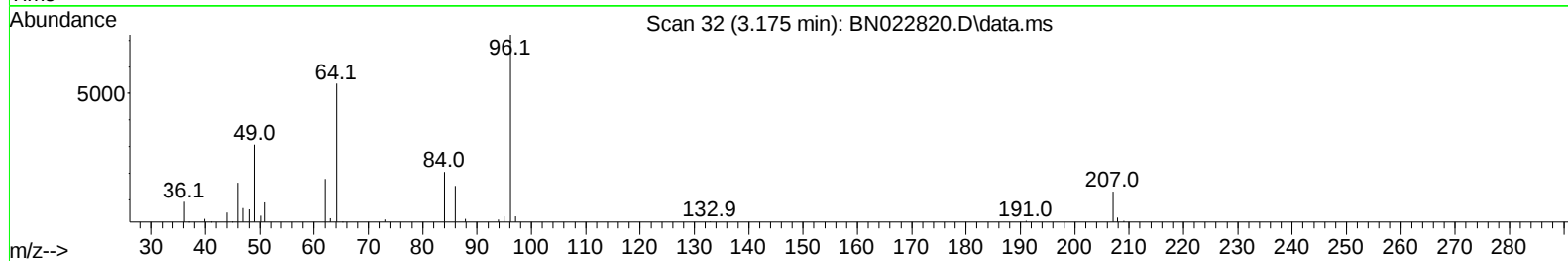
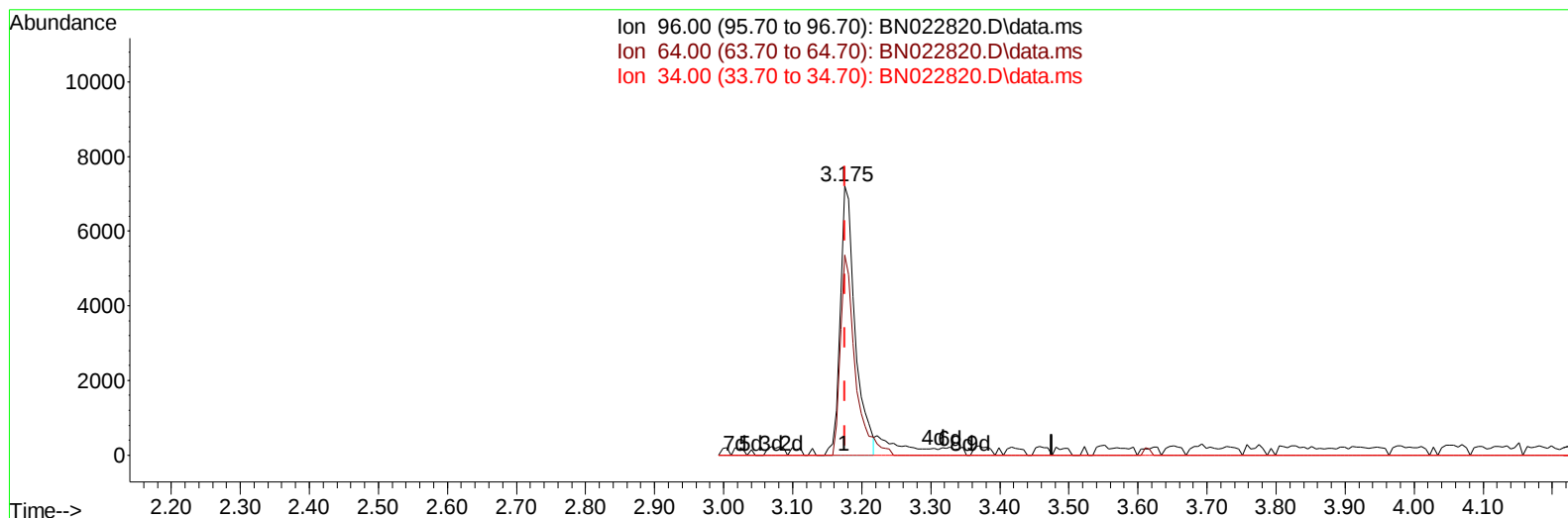
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112122\
 Data File : BN022820.D
 Acq On : 22 Nov 2022 11:56
 Operator : CG/JU
 Sample : PB149042BL
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK042

Manual IntegrationsAPPROVED

Quant Time: Nov 22 12:41:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN11522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 22 05:15:51 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/23/2022
 Supervised By :mohammad ahmed 11/25/2022



TIC: BN022820.D\data.ms

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

3.175min (-0.000) 5.11 ng/uL

response 10888

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	74.35
34.00	0.00	0.00
0.00	0.00	0.00

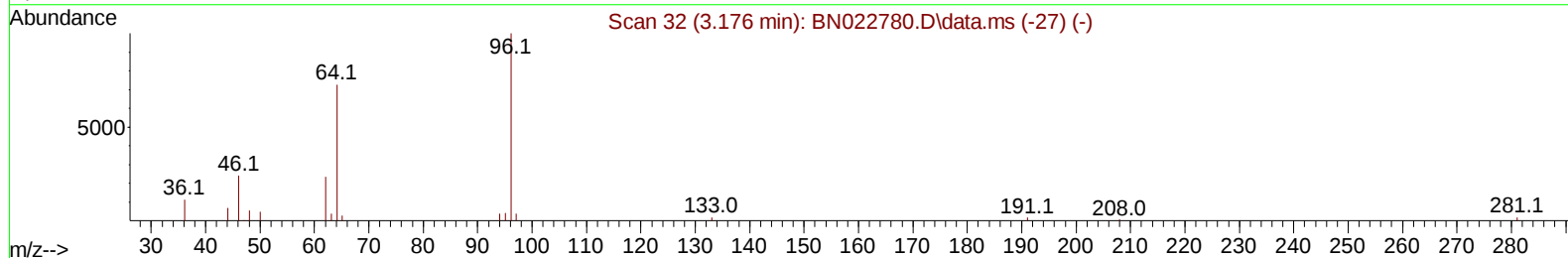
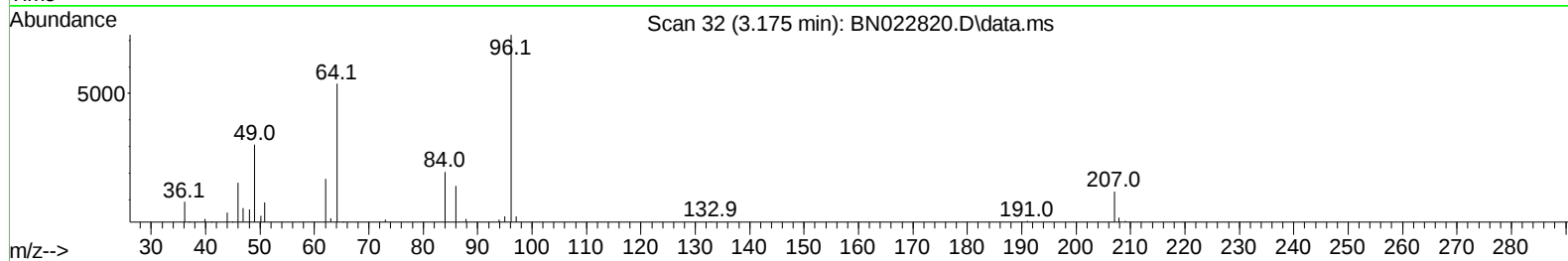
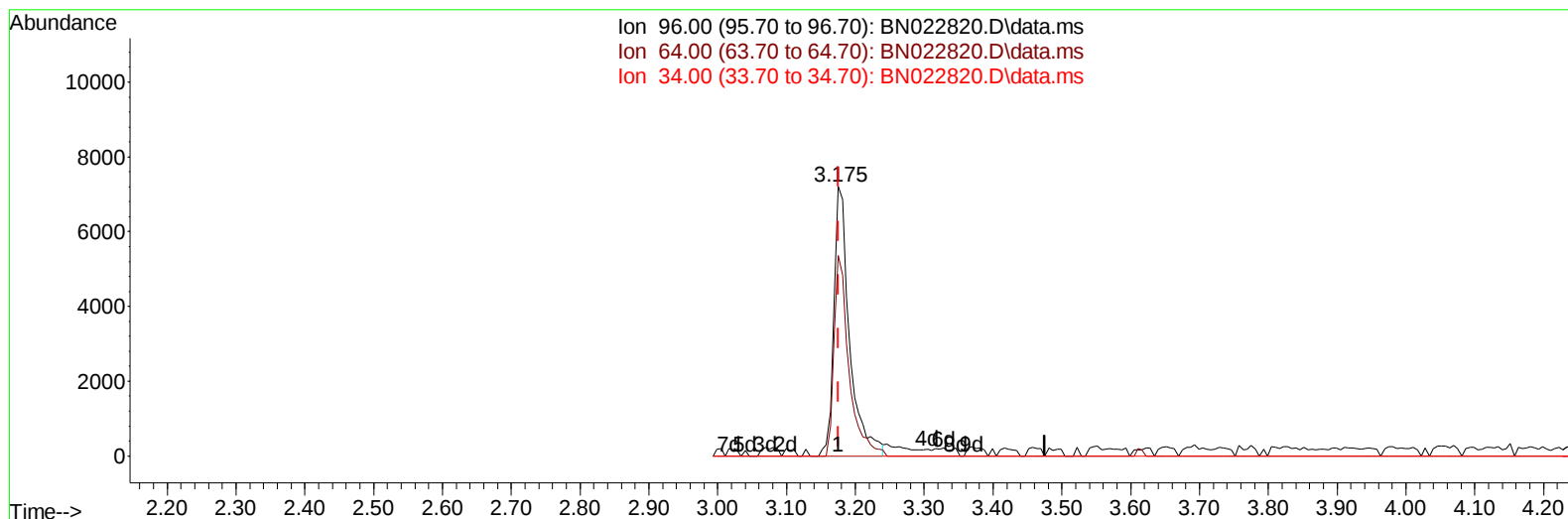
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TIC: BN022820.D\data.ms

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

3.175min (-0.000) 5.37 ng/uL m

response 11455

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	74.35
34.00	0.00	0.00
0.00	0.00	0.00

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 Data File : BN022820.D
 Acq On : 22 Nov 2022 11:56
 Operator : CG/JU
 Sample : PB149042BL
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SBLK042

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Quant Time: Nov 22 12:41:00 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Nov 22 05:15:51 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	72769	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	347222	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.516	164	226167	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.269	188	466740	20.000	ng/ul	0.00
79) Chrysene-d12	21.474	240	448734	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	435065	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.175	96	11455m	5.374	ng/uL	0.00
4) Pyridine-d5	3.611	84	152744	25.341	ng/ul	0.00
7) Phenol-d5	7.010	99	187085	27.221	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	124308	30.166	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	144100	28.629	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	155025	28.571	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	76050	28.296	ng/ul	0.00
24) 2-Nitrophenol-d4	9.751	143	84339	28.123	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	139151	27.218	ng/ul	0.00
31) 4-Chloroaniline-d4	10.810	131	220222	28.279	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	500806	30.443	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	549052	30.291	ng/ul	0.00
54) 4-Nitrophenol-d4	14.734	143	85551	26.087	ng/ul	0.00
60) Fluorene-d10	15.510	176	411805	30.522	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	70841	23.879	ng/ul	0.00
73) Anthracene-d10	17.369	188	658427	32.217	ng/ul	0.00
81) Pyrene-d10	19.674	212	710178	30.539	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	678810	31.038	ng/ul	0.00

Target Compounds Qvalue

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

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