

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112122\
 Data File : BN022827.D
 Acq On : 22 Nov 2022 16:37
 Operator : CG/JU
 Sample : N5593-09
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 H0AD8

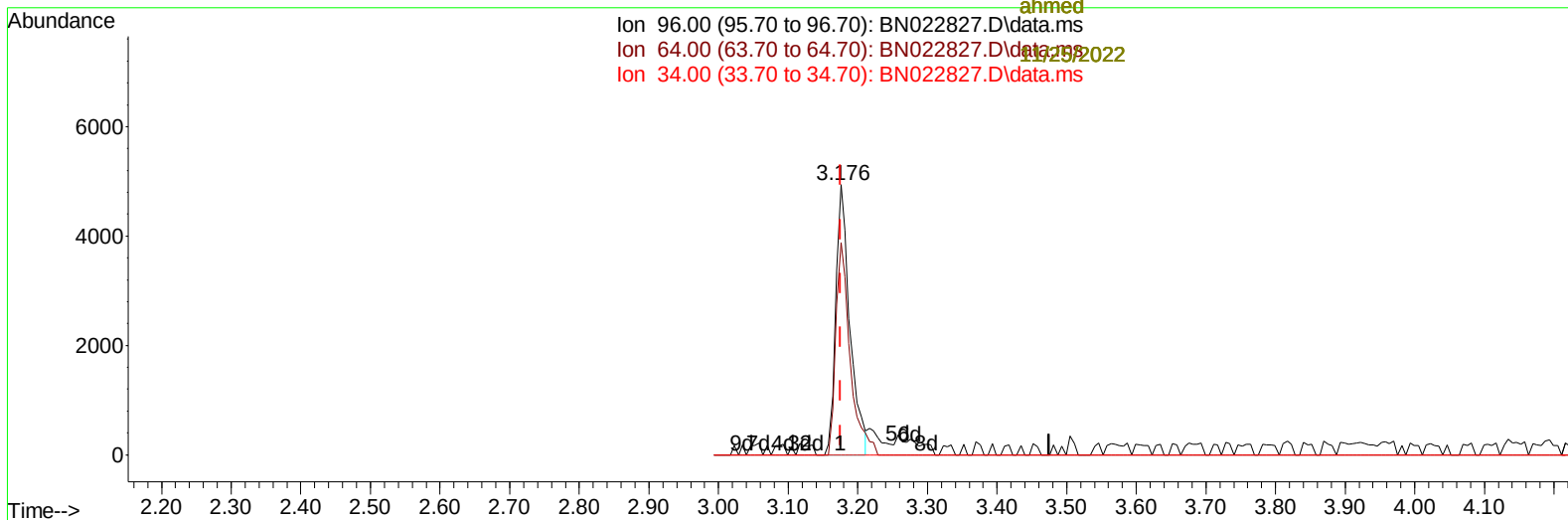
Manual IntegrationsAPPROVED

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

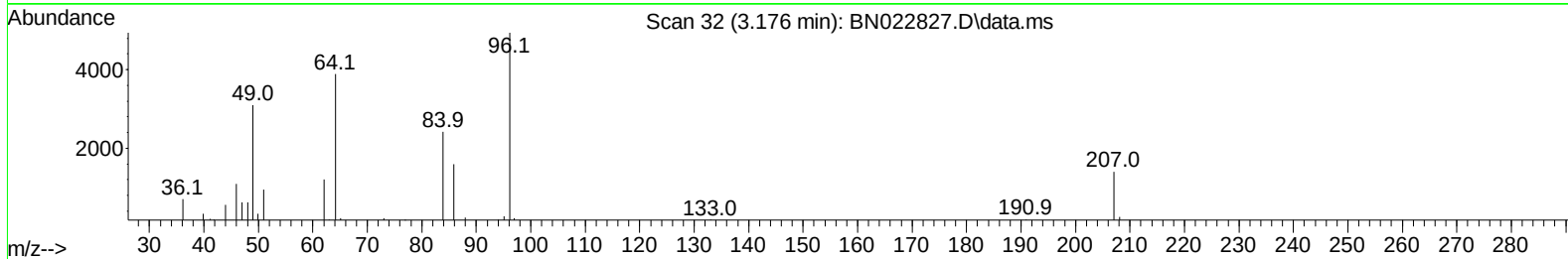
11/23/2022
 Supervised By :mohammad
 ahmed

Quant Time: Nov 22 17:42:55 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

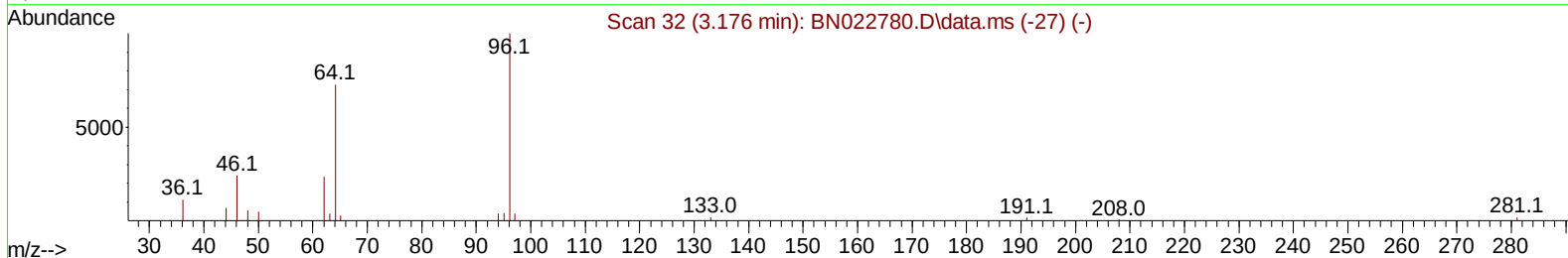
Ion 96.00 (95.70 to 96.70): BN022827.D\data.ms
 Ion 64.00 (63.70 to 64.70): BN022827.D\data.ms
 Ion 34.00 (33.70 to 34.70): BN022827.D\data.ms



Scan 32 (3.176 min): BN022827.D\data.ms



Scan 32 (3.176 min): BN022780.D\data.ms (-27) (-)



TIC: BN022827.D\data.ms

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

3.176min (+ 0.000) 3.82 ng/uL

response 7024

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	78.73
34.00	0.00	0.00
0.00	0.00	0.00

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Quant Time: Nov 22 17:42:55 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)
-----11/25/2022						
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	62798	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	300719	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.516	164	195216	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.269	188	404621	20.000	ng/ul	0.00
79) Chrysene-d12	21.475	240	390158	20.000	ng/ul	0.00
88) Perylene-d12	23.863	264	377920	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.176	96	7534m	4.096	ng/uL	0.00
4) Pyridine-d5	3.617	84	53912	10.364	ng/ul	0.00
7) Phenol-d5	7.011	99	44474	7.498	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	128864	36.237	ng/ul	0.00
11) 2-Chlorophenol-d4	7.369	132	113007	26.017	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	84388	18.022	ng/ul	0.00
21) Nitrobenzene-d5	9.022	128	78733	33.825	ng/ul	0.00
24) 2-Nitrophenol-d4	9.752	143	82029	31.583	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.287	165	130181	29.401	ng/ul	0.00
31) 4-Chloroaniline-d4	10.816	131	184098	27.296	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	540500	38.066	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	577487	36.911	ng/ul	0.00
54) 4-Nitrophenol-d4	14.740	143	20120	7.108	ng/ul	0.00
60) Fluorene-d10	15.510	176	461347	39.615	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	82953	32.254	ng/ul	0.00
73) Anthracene-d10	17.369	188	774607	43.720	ng/ul	0.00
81) Pyrene-d10	19.675	212	869685	43.013	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	824285	43.389	ng/ul	0.00

Target Compounds Qvalue

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

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