

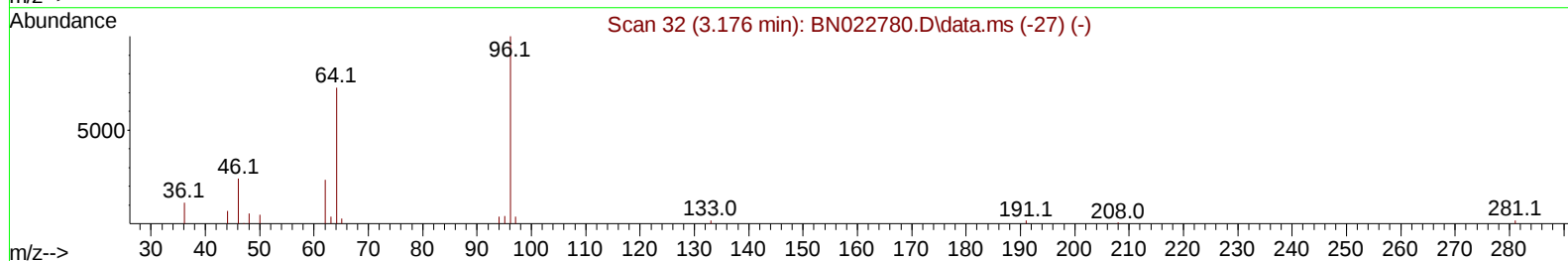
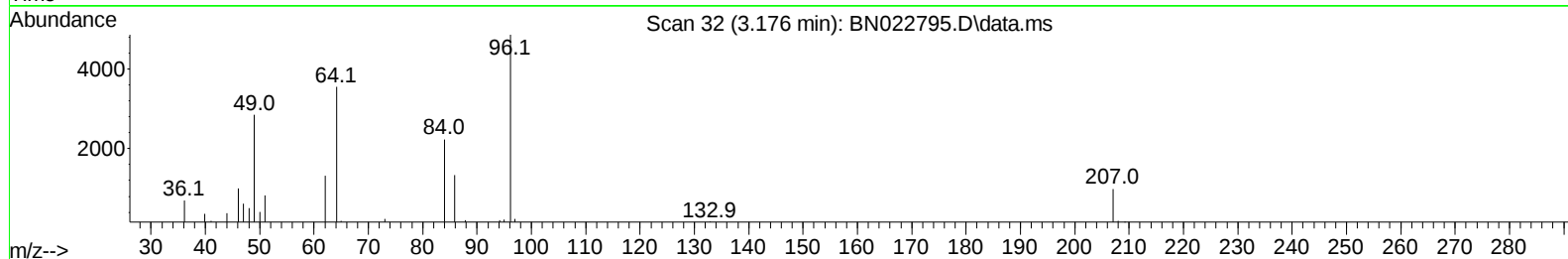
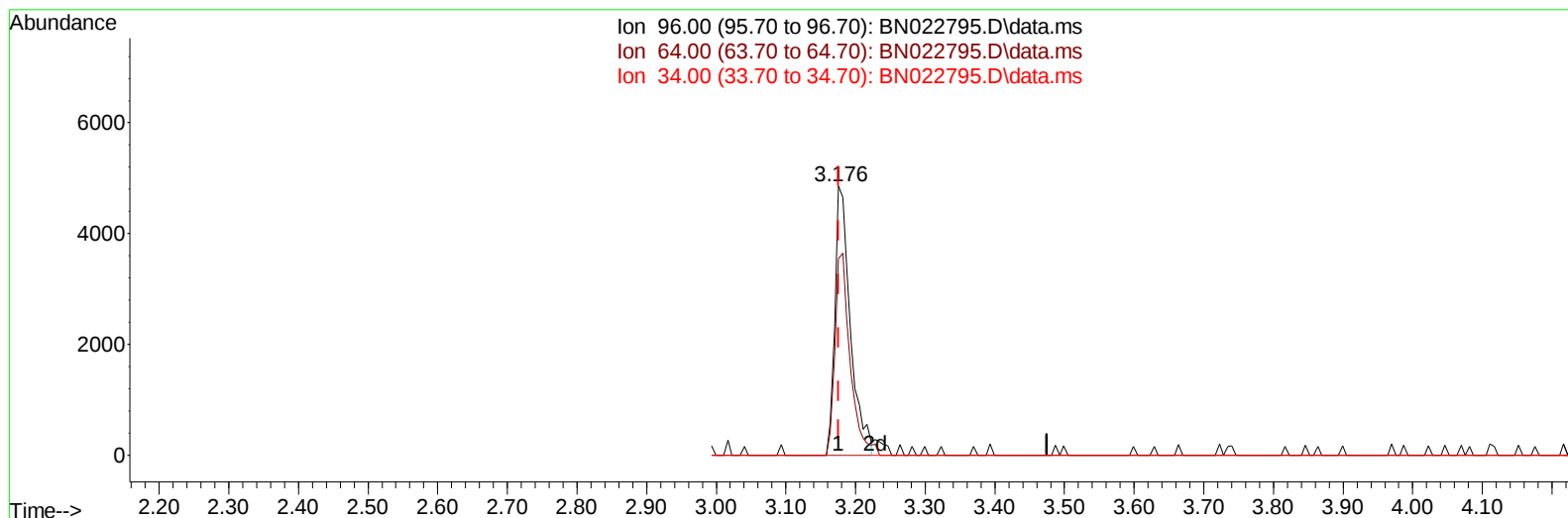
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112122\
Data File : BN022795.D
Acq On : 21 Nov 2022 20:04
Operator : CG/JU
Sample : N5690-13
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGCF4

Manual IntegrationsAPPROVED

Quant Time: Nov 21 21:44:11 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN11522.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Nov 19 05:51:00 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/22/2022
Supervised By :mohammad ahmed 11/25/2022



TIC: BN022795.D\data.ms

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

3.176min (+ 0.000) 3.72 ng/uL

response 7490

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	73.05
34.00	0.00	0.00
0.00	0.00	0.00

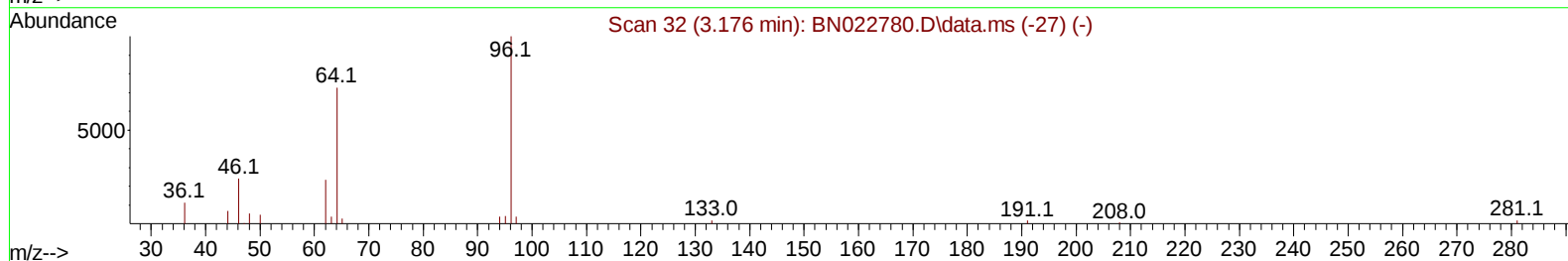
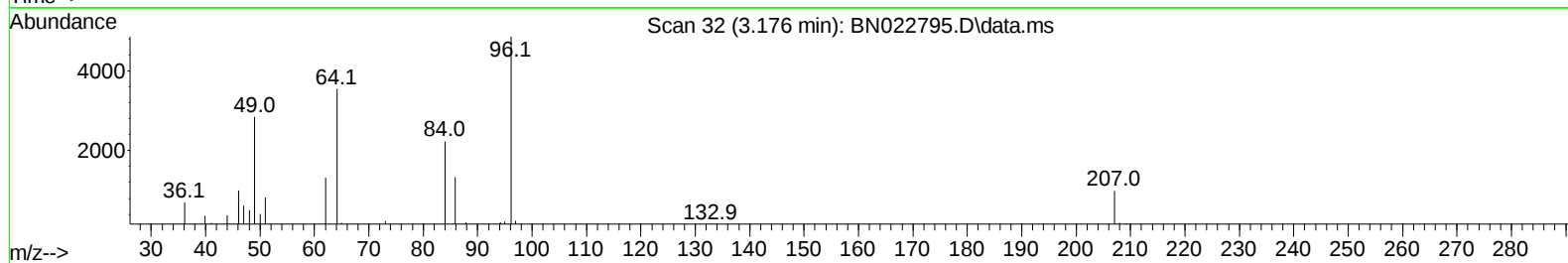
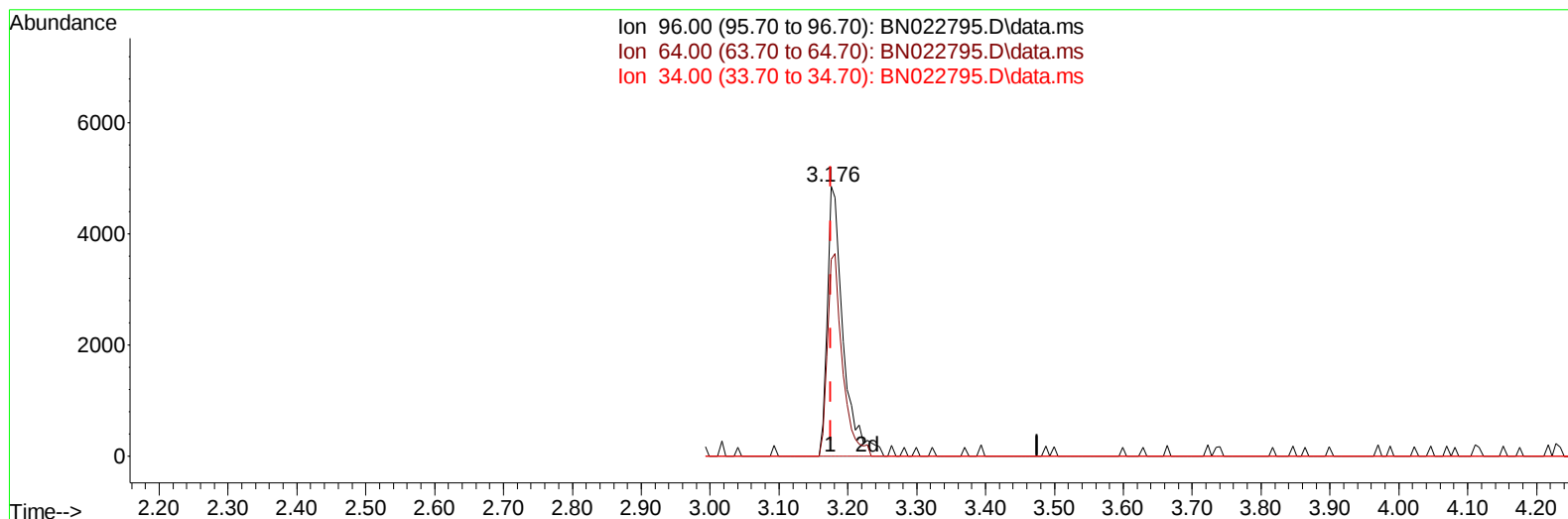
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TIC: BN022795.D\data.ms

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

3.176min (+ 0.000) 3.87 ng/uL m

response 7796

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	73.05
34.00	0.00	0.00
0.00	0.00	0.00

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 Acq On : 21 Nov 2022 20:04
 Operator : CG/JU
 Sample : N5690-13
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
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Manual IntegrationsAPPROVED

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 Supervised By :mohammad ahmed 11/25/2022

Quant Time: Nov 21 21:44:11 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Sat Nov 19 05:51:00 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.846	152	68742	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	330640	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.516	164	219143	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.275	188	450863	20.000	ng/ul	0.00
79) Chrysene-d12	21.475	240	436769	20.000	ng/ul	0.00
88) Perylene-d12	23.869	264	434573	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.176	96	7796m	3.872	ng/uL	0.00
4) Pyridine-d5	3.628	84	19205	3.373	ng/ul	0.02
7) Phenol-d5	7.017	99	39360	6.062	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	121999	31.340	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	104157	21.906	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	75353	14.701	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	76108	29.738	ng/ul	0.00
24) 2-Nitrophenol-d4	9.752	143	79045	27.680	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	121541	24.966	ng/ul	0.00
31) 4-Chloroaniline-d4	10.816	131	177078	23.879	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	521781	32.735	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	572240	32.582	ng/ul	0.00
54) 4-Nitrophenol-d4	14.746	143	16560	5.212	ng/ul	0.00
60) Fluorene-d10	15.516	176	447190	34.207	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	83910	29.280	ng/ul	0.00
73) Anthracene-d10	17.375	188	763606	38.679	ng/ul	0.00
81) Pyrene-d10	19.681	212	862068	38.086	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.716	264	817395	37.417	ng/ul	0.00

Target Compounds Qvalue

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

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