

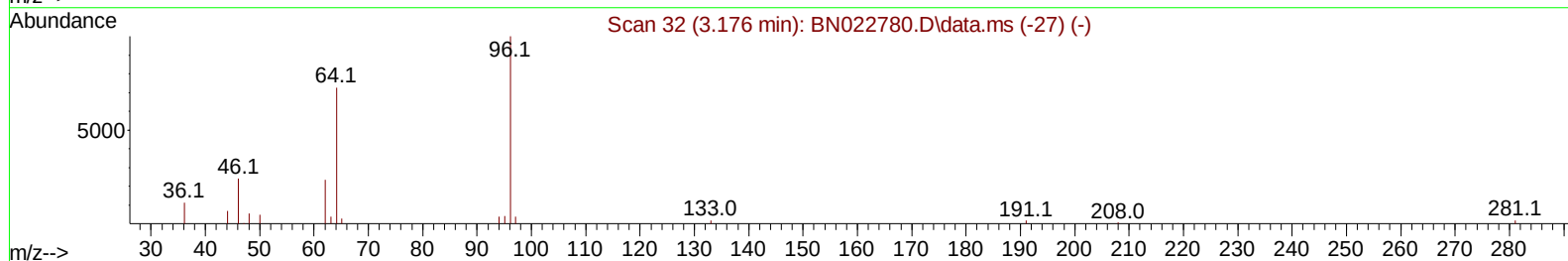
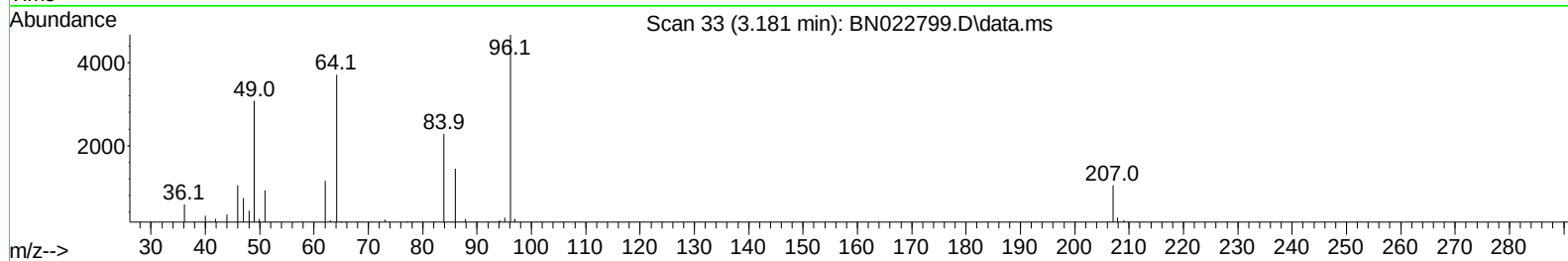
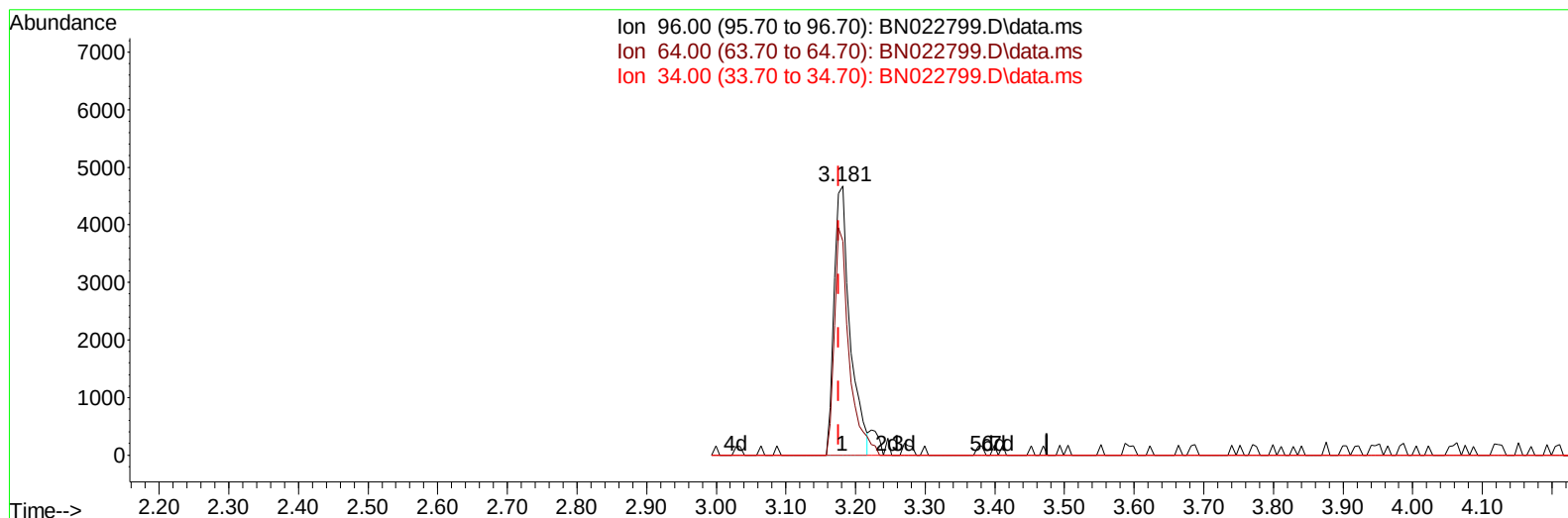
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN112122\
 Data File : BN022799.D
 Acq On : 21 Nov 2022 22:31
 Operator : CG/JU
 Sample : N5690-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGCE6

Manual IntegrationsAPPROVED

Quant Time: Nov 22 00:06:24 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: BN022799.D\data.ms

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

3.181min (+ 0.006) 4.01 ng/uL

response 7438

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	79.47
34.00	0.00	0.00
0.00	0.00	0.00

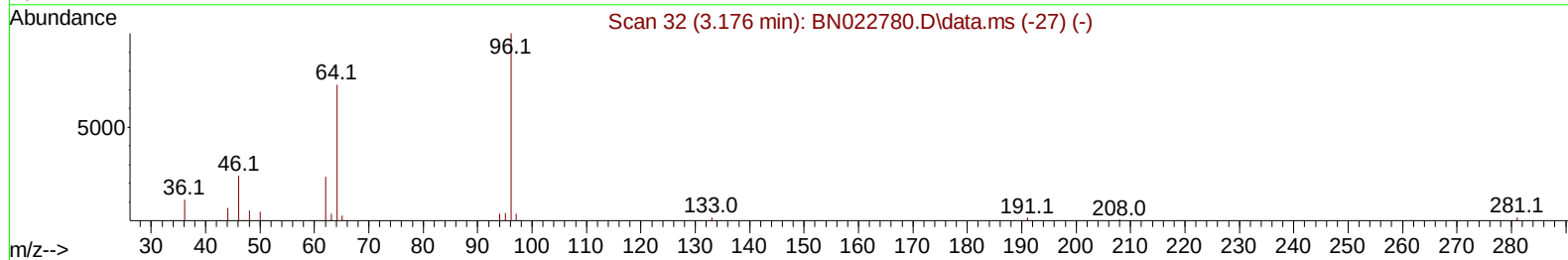
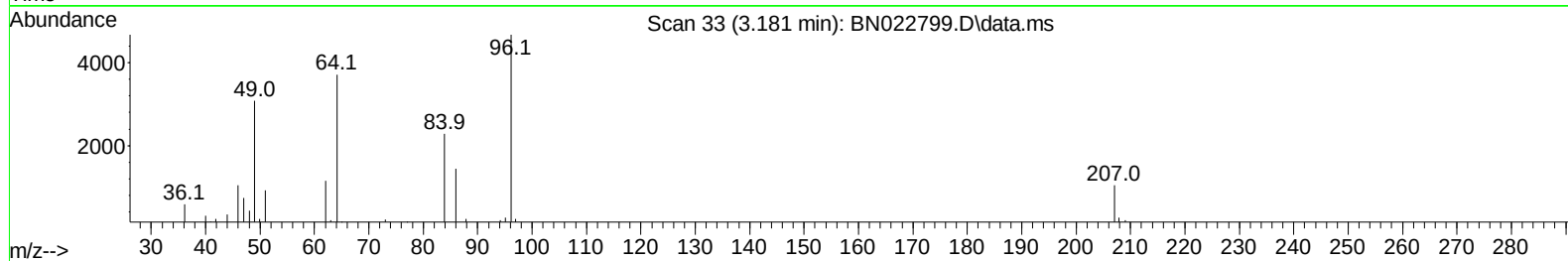
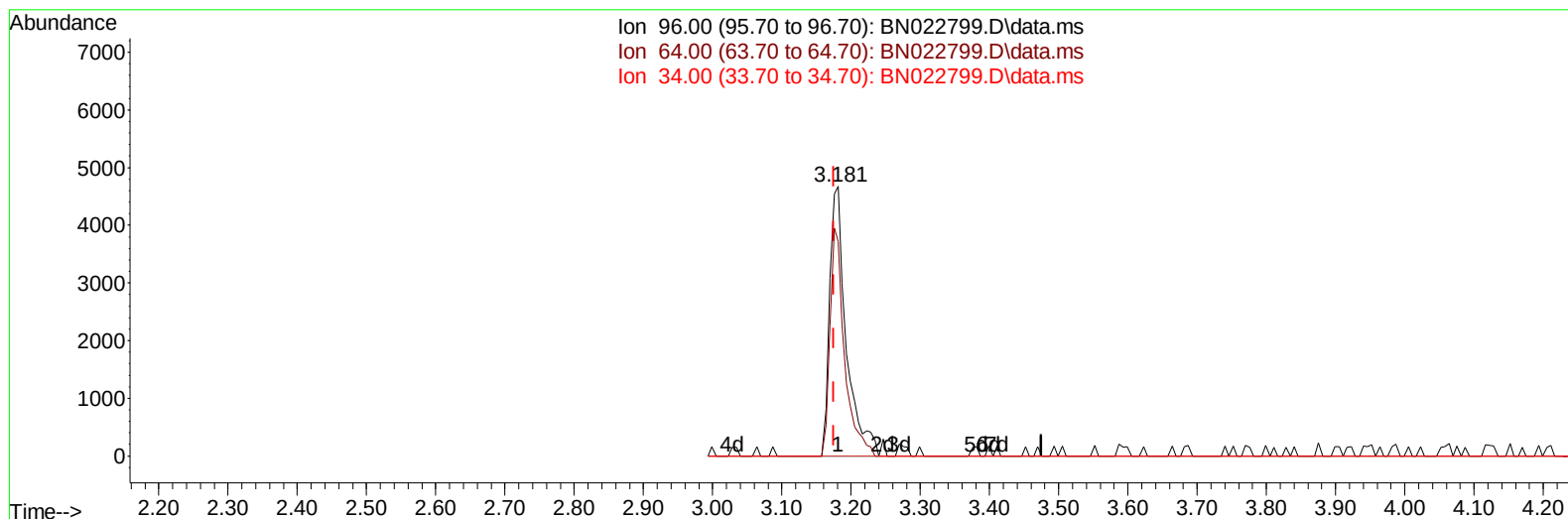
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TIC: BN022799.D\data.ms

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

3.181min (+ 0.006) 4.23 ng/uL m

response 7833

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	79.47
34.00	0.00	0.00
0.00	0.00	0.00

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 Acq On : 21 Nov 2022 22:31
 Operator : CG/JU
 Sample : N5690-02
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
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Quant Time: Nov 22 00:06:24 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	63256	20.000	ng/ul	0.00
20) Naphthalene-d8	10.663	136	299715	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.516	164	197562	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.275	188	416239	20.000	ng/ul	0.00
79) Chrysene-d12	21.475	240	406705	20.000	ng/ul	0.00
88) Perylene-d12	23.869	264	394771	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	7833m	4.228	ng/uL	0.00
4) Pyridine-d5	3.623	84	46831	8.938	ng/ul	0.01
7) Phenol-d5	7.017	99	44738	7.488	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.181	67	134835	37.642	ng/ul	0.00
11) 2-Chlorophenol-d4	7.375	132	115371	26.369	ng/ul	0.00
15) 4-Methylphenol-d8	8.569	113	85364	18.099	ng/ul	0.00
21) Nitrobenzene-d5	9.028	128	84848	36.574	ng/ul	0.00
24) 2-Nitrophenol-d4	9.752	143	87634	33.854	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.293	165	135361	30.673	ng/ul	0.00
31) 4-Chloroaniline-d4	10.816	131	192490	28.636	ng/ul	0.00
46) Dimethylphthalate-d6	13.934	166	573529	39.912	ng/ul	0.00
49) Acenaphthylene-d8	14.210	160	612981	38.715	ng/ul	0.00
54) 4-Nitrophenol-d4	14.746	143	17533	6.120	ng/ul	0.00
60) Fluorene-d10	15.516	176	477652	40.529	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	81227	30.702	ng/ul	0.00
73) Anthracene-d10	17.375	188	764846	41.964	ng/ul	0.00
81) Pyrene-d10	19.675	212	872120	41.378	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.710	264	842278	42.444	ng/ul	-0.01

Target Compounds Qvalue

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

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