

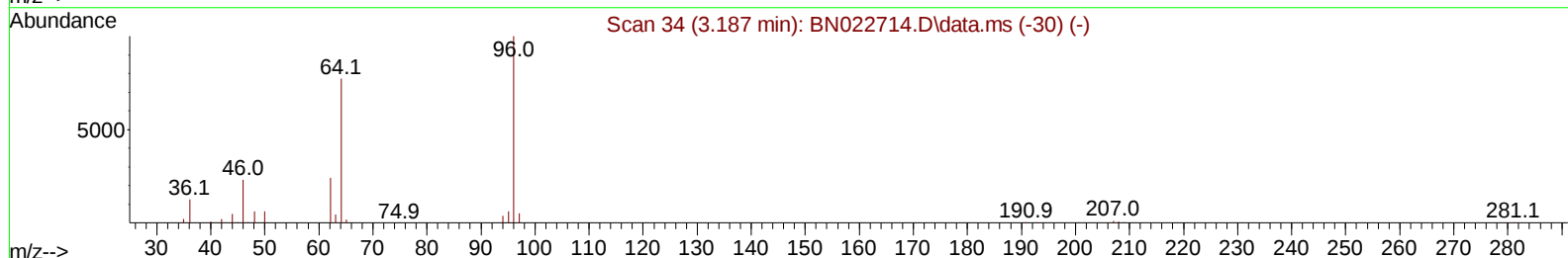
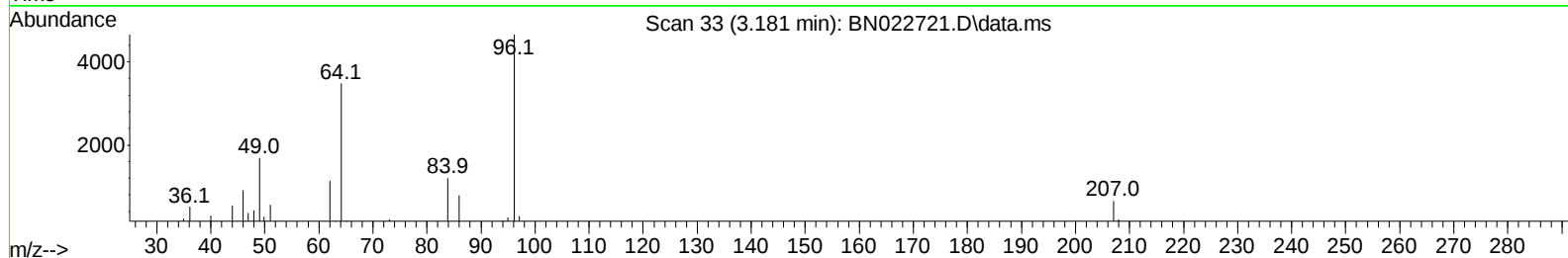
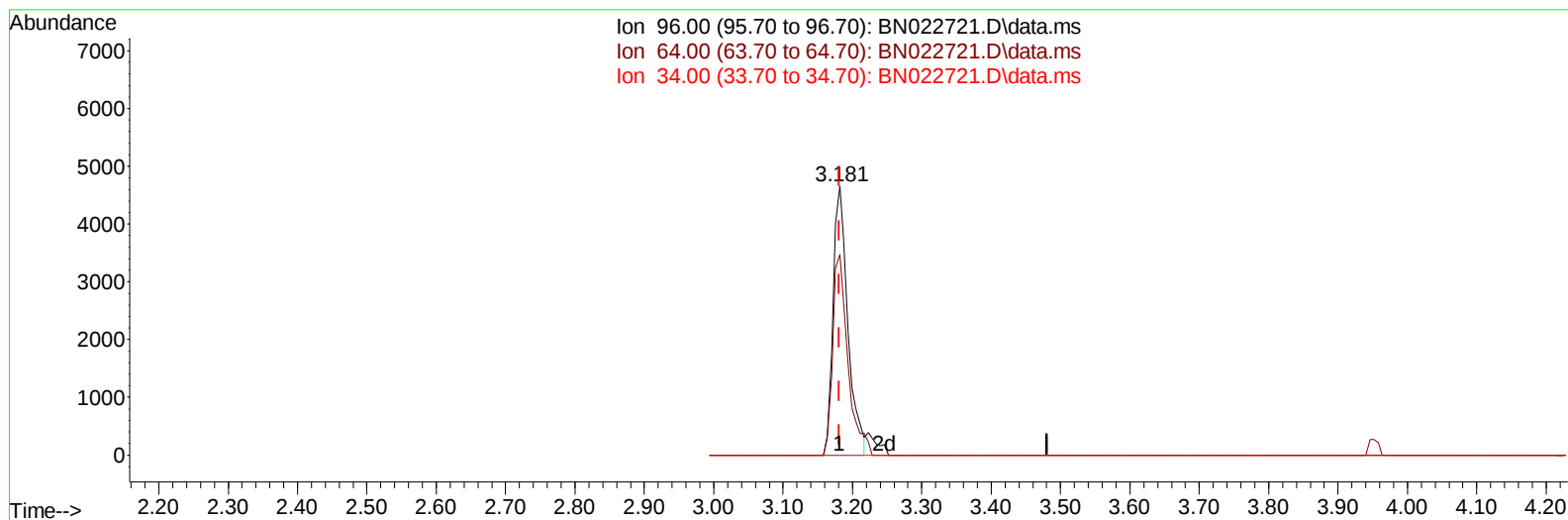
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111722\
 Data File : BN022721.D
 Acq On : 17 Nov 2022 15:51
 Operator : CG/JU
 Sample : N5622-15
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BGCA0

Manual IntegrationsAPPROVED

Quant Time: Nov 17 22:11:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN111522.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Nov 17 22:08:43 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 11/18/2022
 Supervised By :mohammad ahmed 11/21/2022



TIC: BN022721.D\data.ms

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

3.181min (+ 0.000) 5.07 ng/uL

response 6839

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	74.68
34.00	0.00	0.00
0.00	0.00	0.00

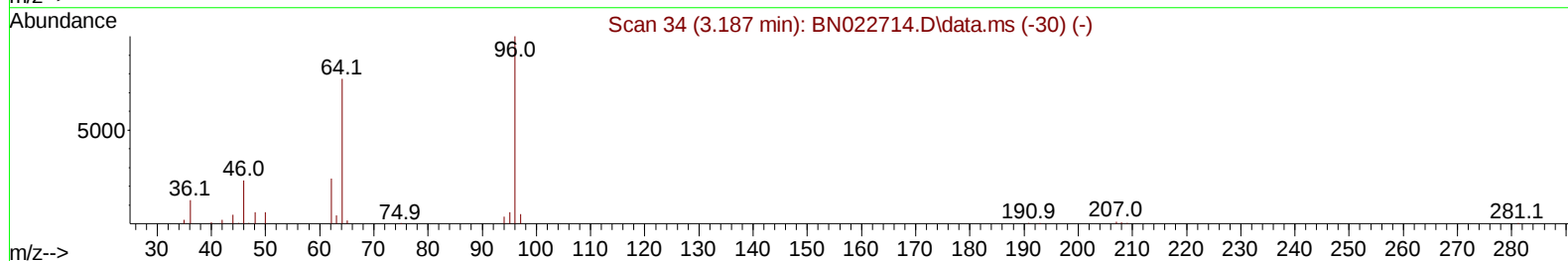
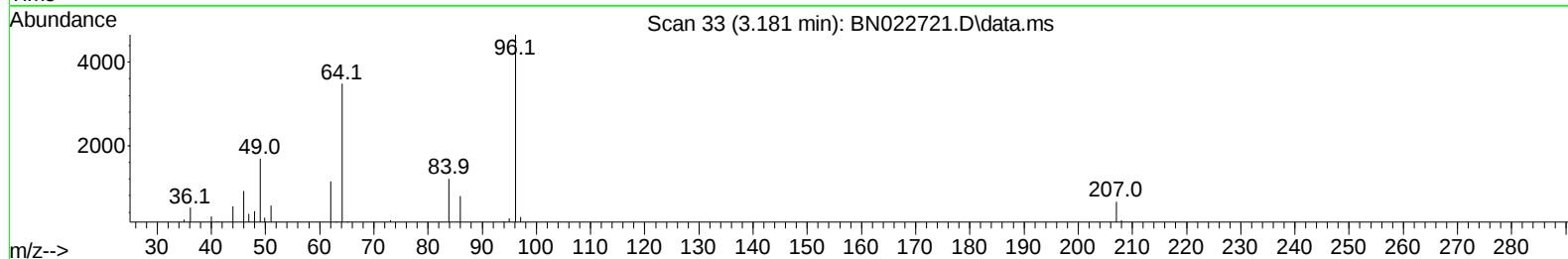
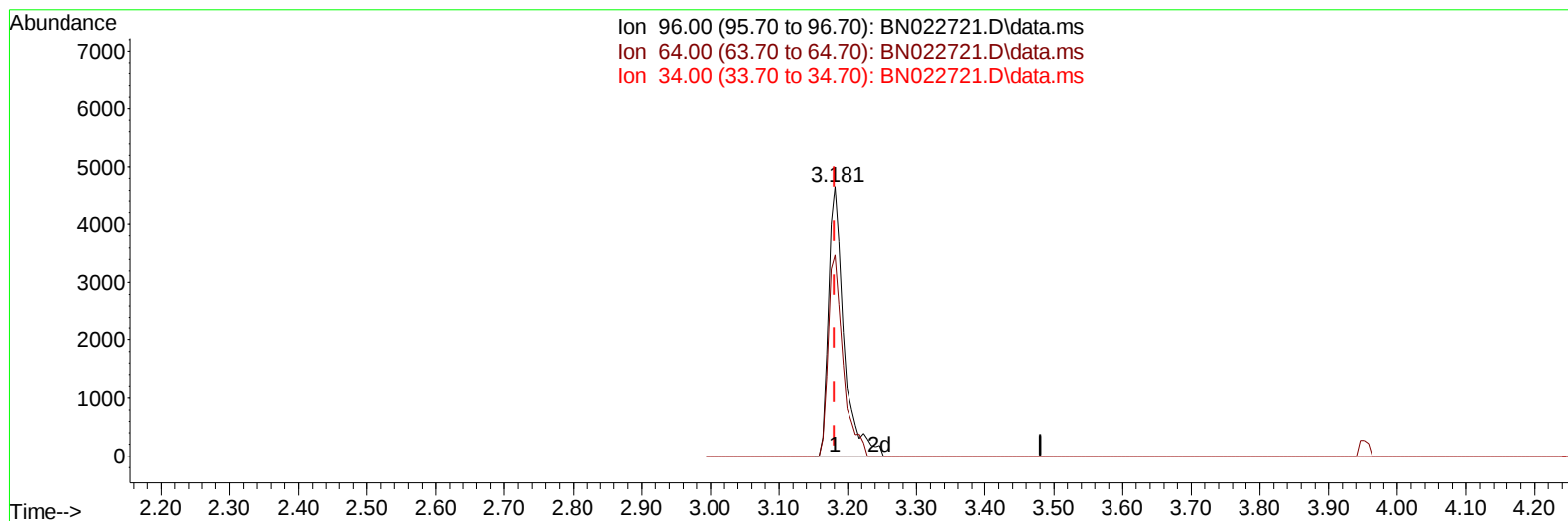
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(3) 1,4-Dioxane-d8 (S)

3.181min (+ 0.000) 5.39 ng/uL m

response 7266

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.50	74.68
34.00	0.00	0.00
0.00	0.00	0.00

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.858	152	46060	20.000	ng/ul	0.00
20) Naphthalene-d8	10.675	136	225966	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.528	164	156077	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.287	188	341270	20.000	ng/ul	0.00
79) Chrysene-d12	21.492	240	353378	20.000	ng/ul	0.00
88) Perylene-d12	23.892	264	369881	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.181	96	7266m	5.386	ng/uL	0.00
4) Pyridine-d5	3.628	84	31008	8.127	ng/ul	0.01
7) Phenol-d5	7.028	99	35707	8.208	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.187	67	105342	40.387	ng/ul	0.00
11) 2-Chlorophenol-d4	7.387	132	93172	29.245	ng/ul	0.00
15) 4-Methylphenol-d8	8.581	113	66982	19.503	ng/ul	0.00
21) Nitrobenzene-d5	9.040	128	64195	36.702	ng/ul	0.00
24) 2-Nitrophenol-d4	9.763	143	69004	35.357	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.305	165	103852	31.214	ng/ul	0.00
31) 4-Chloroaniline-d4	10.828	131	171766	33.893	ng/ul	0.00
46) Dimethylphthalate-d6	13.946	166	460251	40.542	ng/ul	0.00
49) Acenaphthylene-d8	14.222	160	478079	38.220	ng/ul	0.00
54) 4-Nitrophenol-d4	14.757	143	15160	6.699	ng/ul	0.00
60) Fluorene-d10	15.528	176	382199	41.049	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.657	200	71992	33.189	ng/ul	0.00
73) Anthracene-d10	17.387	188	662542	44.337	ng/ul	0.00
81) Pyrene-d10	19.692	212	748177	40.854	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.745	264	829133	44.593	ng/ul	0.00

Target Compounds Qvalue

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

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