

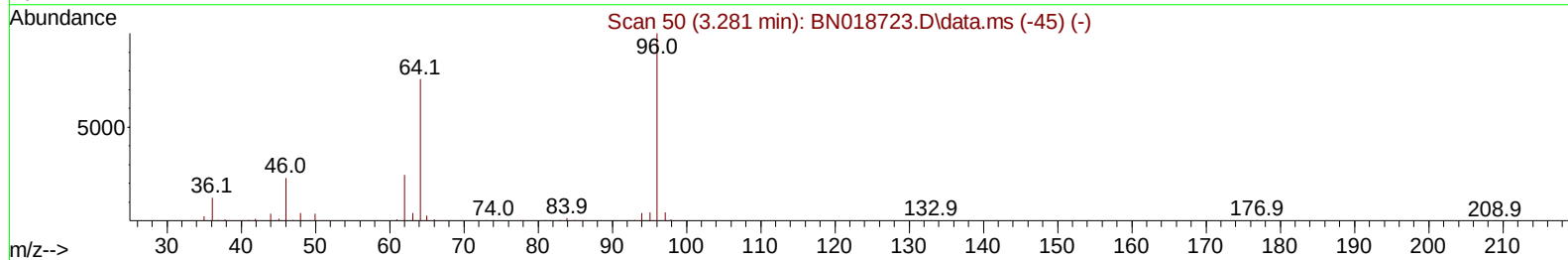
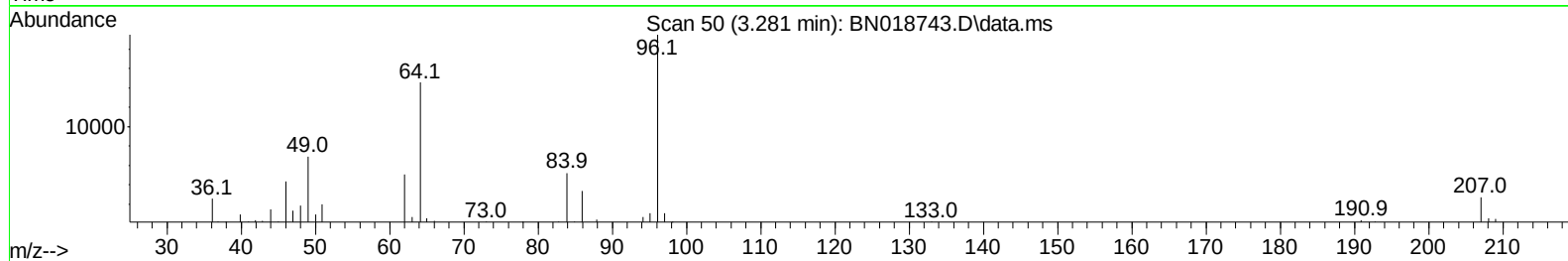
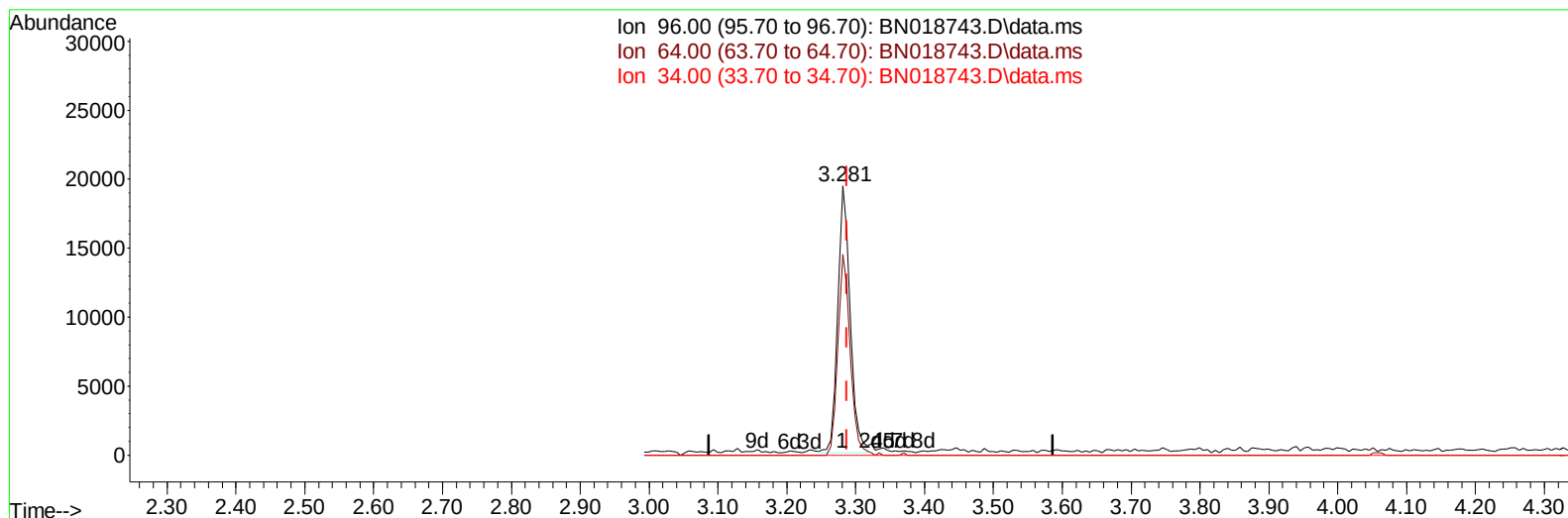
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN022522\
Data File : BN018743.D
Acq On : 25 Feb 2022 23:03
Operator : CG/JU
Sample : N1656-13
Misc :
ALS Vial : 22 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
BGQ20

Manual IntegrationsAPPROVED

Quant Time: Feb 26 01:13:52 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN021622.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Feb 16 16:48:14 2022
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 02/28/2022
Supervised By :mohammad ahmed 02/28/2022



TIC: BN018743.D\data.ms

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

3.281min (-0.006) 4.72 ng/uL

response 24399

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.80	74.69
34.00	0.00	0.00
0.00	0.00	0.00

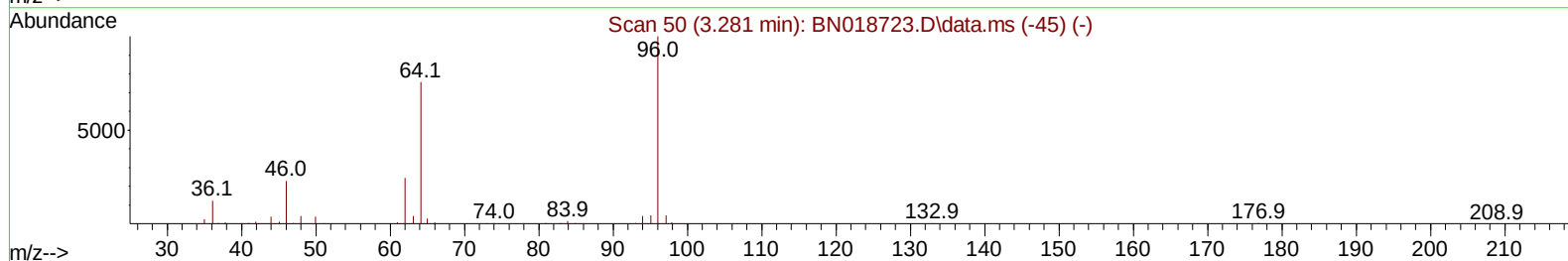
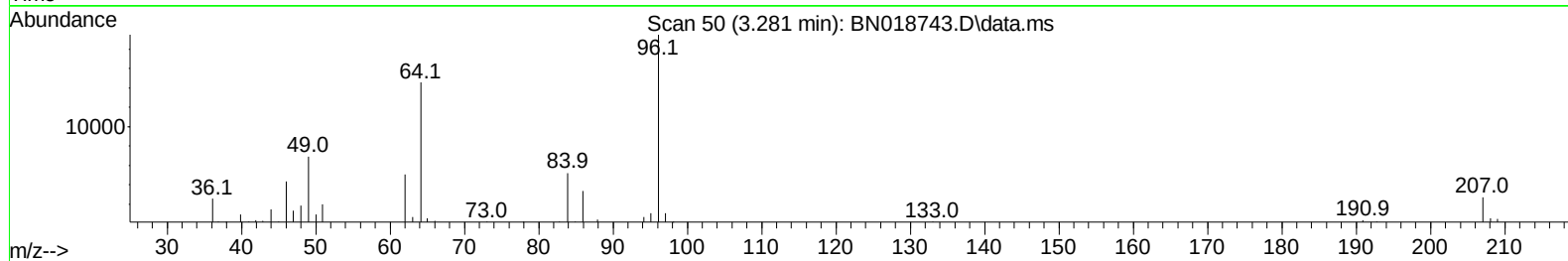
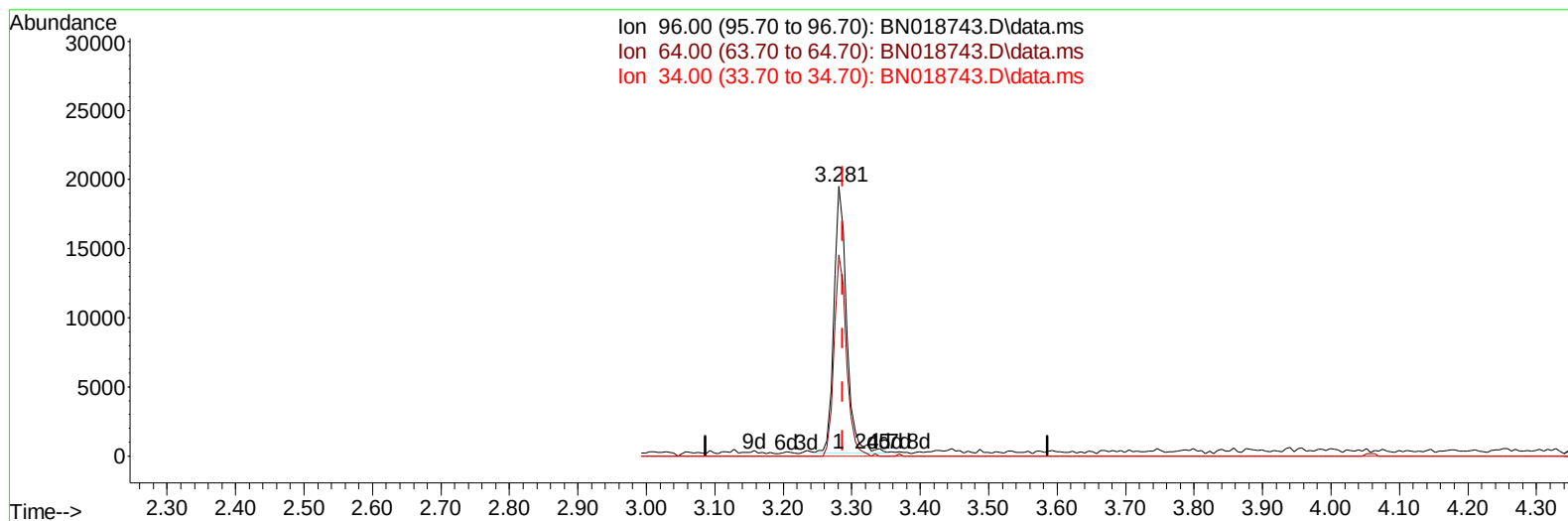
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TIC: BN018743.D\data.ms

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

3.281min (-0.006) 4.80 ng/uL m

response 24800

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	73.80	74.69
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	7.910	152	194897	20.000	ng/ul	0.00
20) Naphthalene-d8	10.716	136	731599	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.545	164	407505	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.280	188	811487	20.000	ng/ul	0.00
79) Chrysene-d12	21.463	240	825129	20.000	ng/ul	0.00
88) Perylene-d12	23.857	264	878744	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.281	96	24800m	4.796	ng/uL	0.00
4) Pyridine-d5	3.705	84	73062	5.226	ng/ul	0.00
7) Phenol-d5	7.057	99	106930	6.652	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.228	67	285225	28.590	ng/ul	0.00
11) 2-Chlorophenol-d4	7.434	132	294827	23.288	ng/ul	0.00
15) 4-Methylphenol-d8	8.610	113	181309	14.770	ng/ul	0.00
21) Nitrobenzene-d5	9.069	128	158958	28.276	ng/ul	0.00
24) 2-Nitrophenol-d4	9.793	143	167052	27.746	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.328	165	291272	25.346	ng/ul	-0.01
31) 4-Chloroaniline-d4	10.846	131	395256	25.742	ng/ul	0.00
46) Dimethylphthalate-d6	13.945	166	895152	30.344	ng/ul	0.00
49) Acenaphthylene-d8	14.239	160	1146647	30.002	ng/ul	0.00
54) 4-Nitrophenol-d4	14.722	143	24609	4.801	ng/ul	0.00
60) Fluorene-d10	15.534	176	802235	31.499	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.645	200	109226	22.813	ng/ul	0.00
73) Anthracene-d10	17.380	188	1302151	33.535	ng/ul	0.00
81) Pyrene-d10	19.669	212	1551226	32.119	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.704	264	1581350	35.286	ng/ul	0.00

Target Compounds Qvalue

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

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