

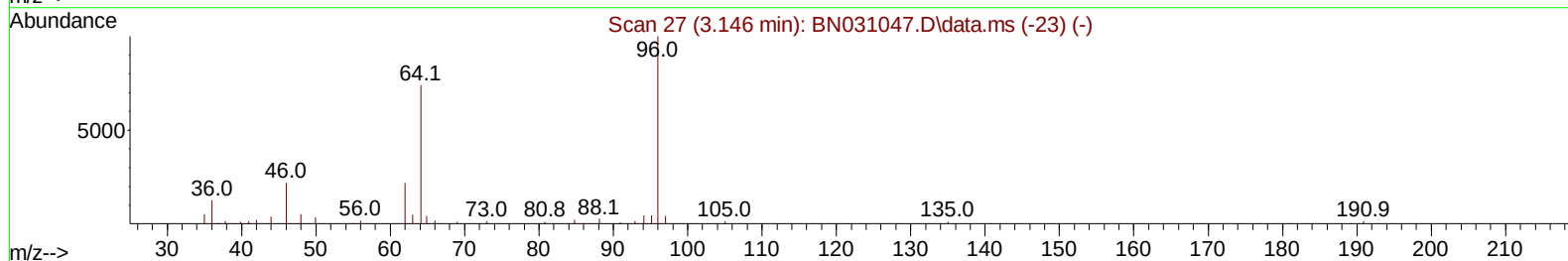
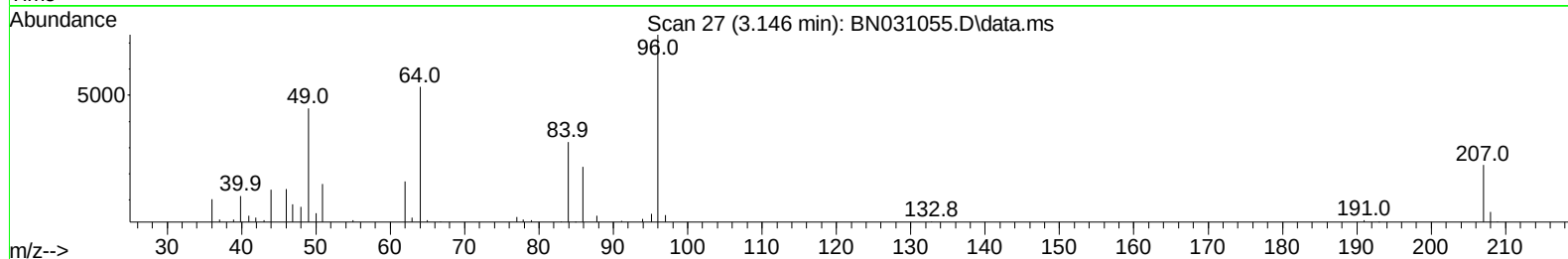
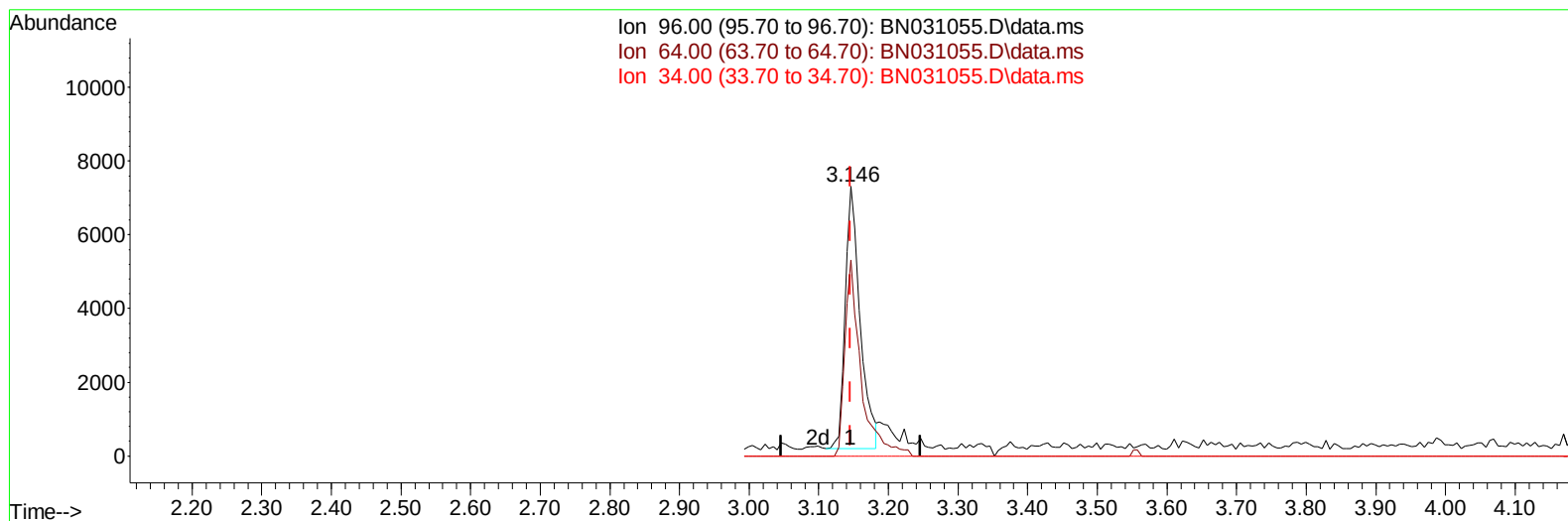
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041924\
Data File : BN031055.D
Acq On : 20 Apr 2024 01:40
Operator : MA/JU
Sample : PB160347BL
Misc :
ALS Vial : 3 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
SBLK347

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/22/2024
Supervised By :mohammad ahmed 04/22/2024

Quant Time: Apr 20 02:09:47 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN041924.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Apr 19 22:56:25 2024
Response via : Initial Calibration



TIC: BN031055.D\data.ms

(3) 1,4-Dioxane-d8 (S)**3.146min (+ 0.000) 3.98 ng/uL****response 10689**

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.80	72.75
34.00	0.00	0.00
0.00	0.00	0.00

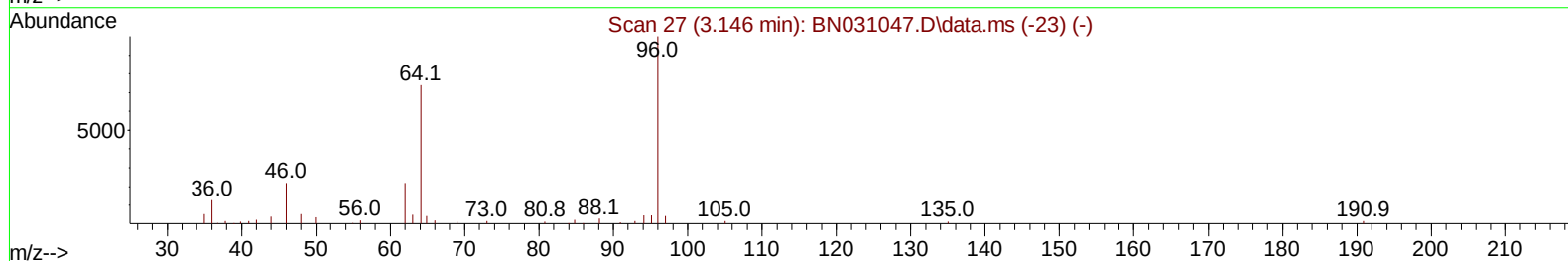
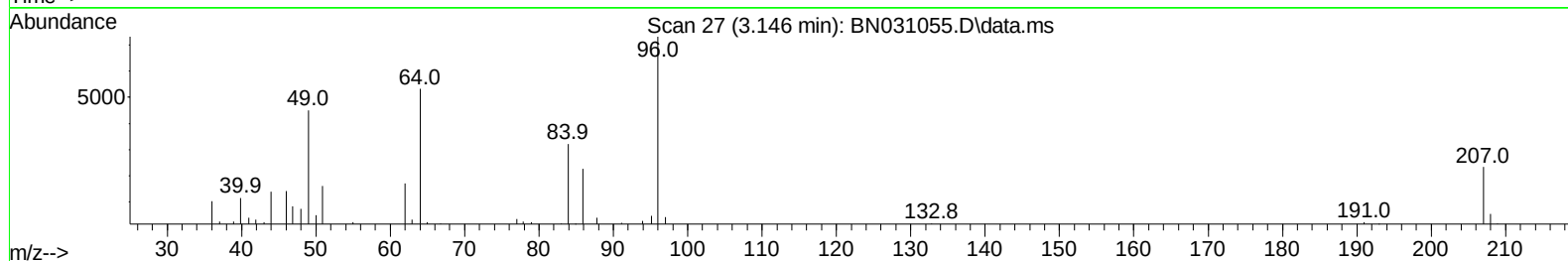
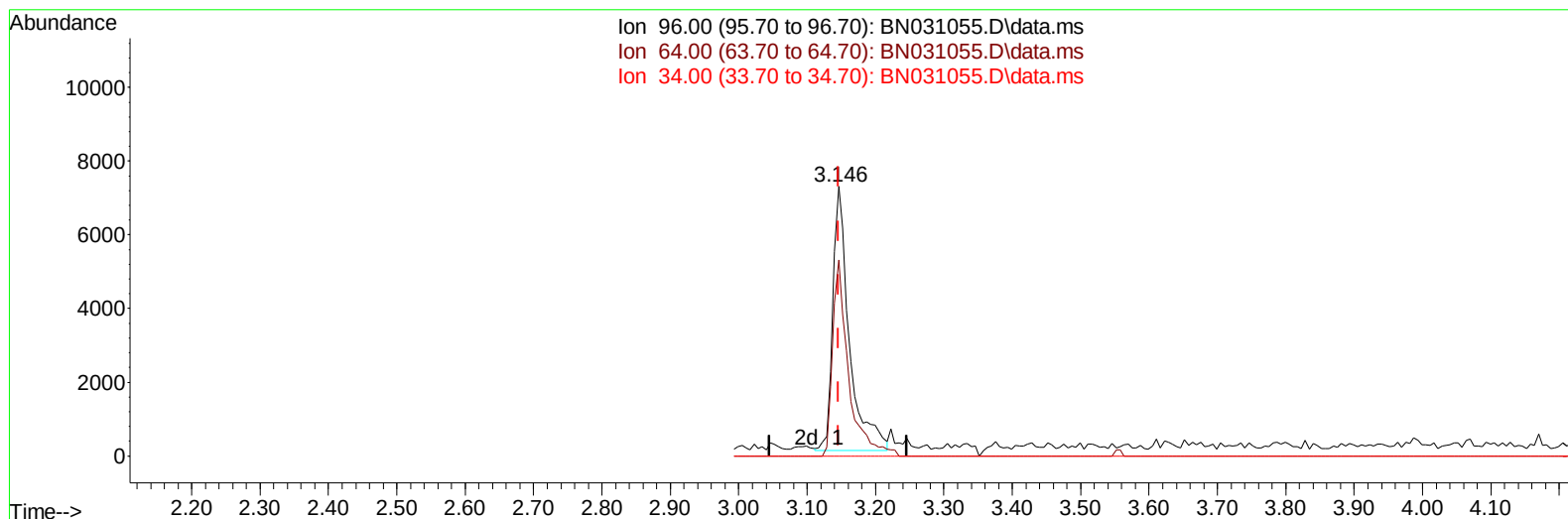
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TIC: BN031055.D\data.ms

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

3.146min (+ 0.000) 4.50 ng/uL m

response 12072

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	72.80	72.75
34.00	0.00	0.00
0.00	0.00	0.00

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 QLast Update : Fri Apr 19 22:56:25 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.634	152	125445	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	579364	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.281	164	403588	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.034	188	919710	20.000	ng/ul	0.00
79) Chrysene-d12	21.263	240	1092181	20.000	ng/ul	0.00
88) Perylene-d12	24.169	264	1489394	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	12072m	4.496	ng/uL	0.00
4) Pyridine-d5	3.552	84	134152	17.349	ng/ul	0.00
7) Phenol-d5	6.811	99	223388	21.110	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.975	67	135129	21.948	ng/ul	0.00
11) 2-Chlorophenol-d4	7.169	132	179523	21.788	ng/ul	0.00
15) 4-Methylphenol-d8	8.340	113	186958	21.008	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	94785	22.770	ng/ul	0.00
24) 2-Nitrophenol-d4	9.505	143	96332	22.263	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.040	165	178769	20.950	ng/ul	0.00
31) 4-Chloroaniline-d4	10.557	131	244824	18.360	ng/ul	0.00
46) Dimethylphthalate-d6	13.698	166	733483	26.117	ng/ul	0.00
49) Acenaphthylene-d8	13.975	160	790831	24.849	ng/ul	0.00
54) 4-Nitrophenol-d4	14.493	143	127903	21.820	ng/ul	0.00
60) Fluorene-d10	15.281	176	629072	25.730	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	89735	19.339	ng/ul	0.00
73) Anthracene-d10	17.128	188	1034421	25.949	ng/ul	0.00
81) Pyrene-d10	19.433	212	1320453	24.604	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.969	264	1807793	25.596	ng/ul	0.00

Target Compounds Qvalue

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

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