

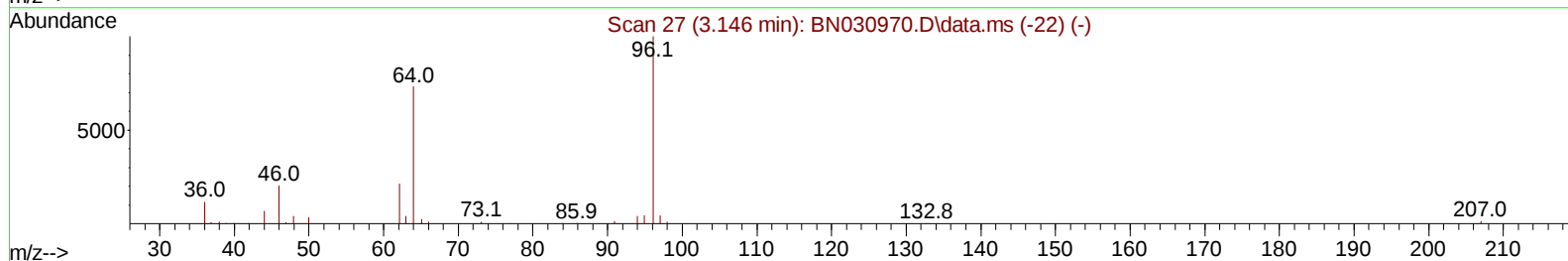
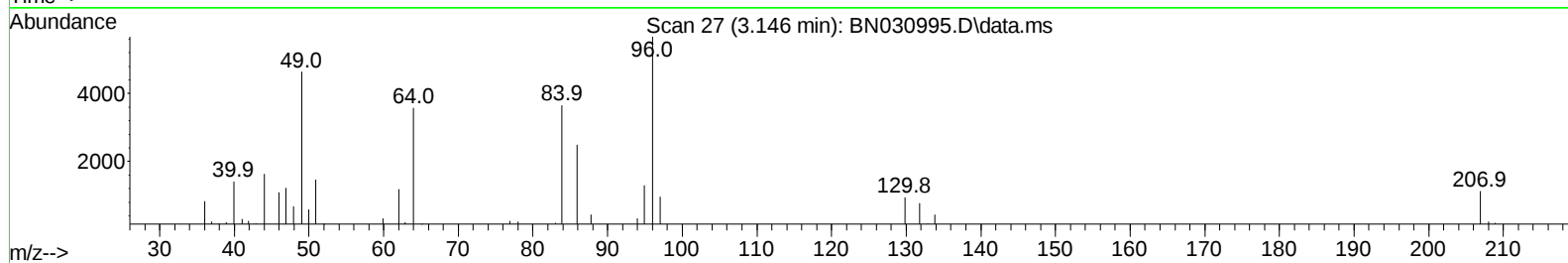
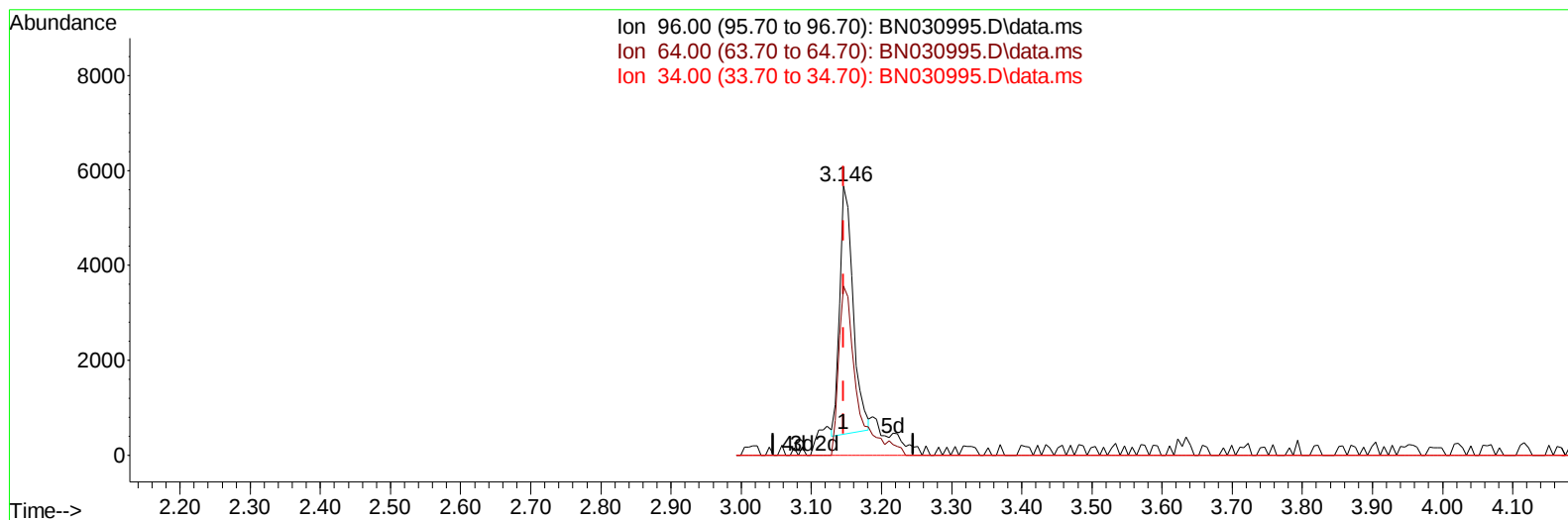
Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN041724\
Data File : BN030995.D
Acq On : 18 Apr 2024 04:00
Operator : MA/JU
Sample : P2053-03
Misc :
ALS Vial : 28 Sample Multiplier: 1

Instrument :
BNA_N
ClientSampleId :
YC558

Manual IntegrationsAPPROVED

Quant Time: Apr 18 04:30:14 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Apr 17 22:24:16 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 04/19/2024
Supervised By :mohammad ahmed 04/30/2024



TIC: BN030995.D\data.ms

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

3.146min (+ 0.000) 3.07 ng/uL

response 6968

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	65.50	62.95
34.00	0.00	0.00
0.00	0.00	0.00

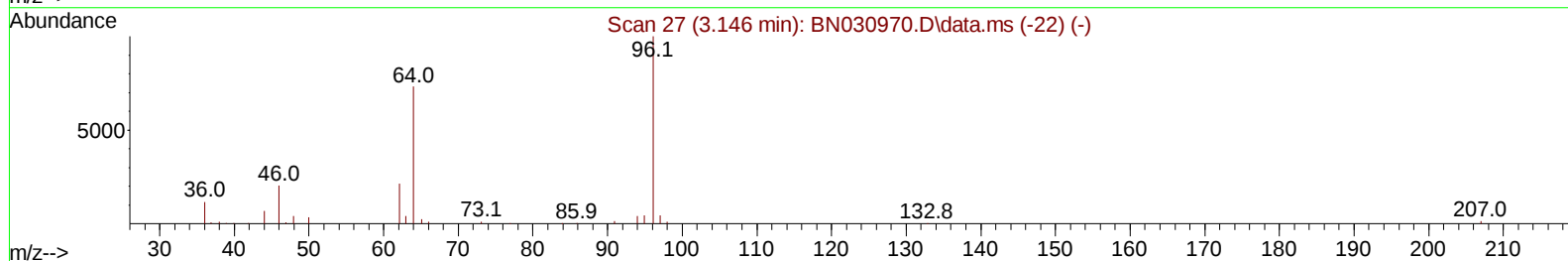
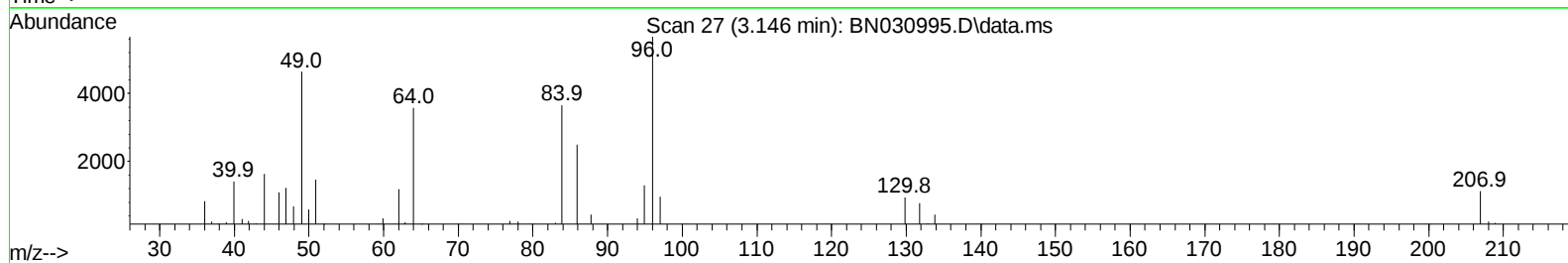
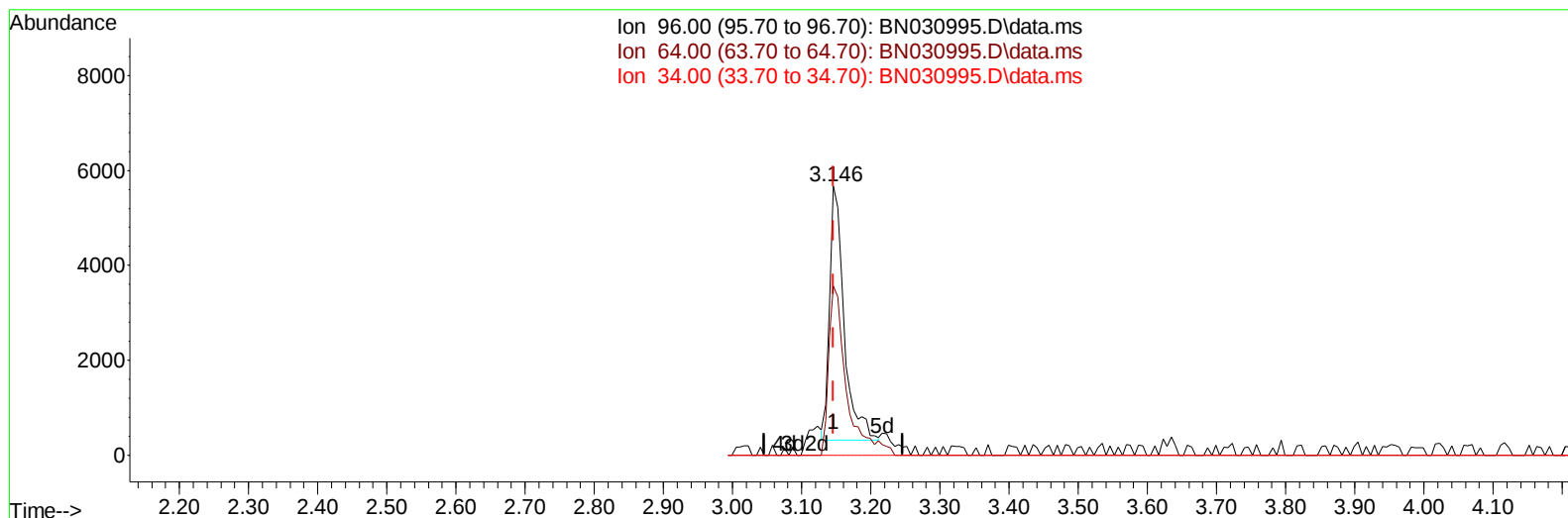
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TIC: BN030995.D\data.ms

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

3.146min (+ 0.000) 3.46 ng/uL m

response 7863

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	65.50	62.95
34.00	0.00	0.00
0.00	0.00	0.00

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 Sample : P2053-03
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 YCSS8

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 04/19/2024
 Supervised By :mohammad ahmed 04/30/2024

Quant Time: Apr 18 04:31:16 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-BN032824.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Apr 17 22:24:16 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.640	152	106530	20.000	ng/ul	0.00
20) Naphthalene-d8	10.416	136	496774	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.287	164	352855	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.034	188	807280	20.000	ng/ul	0.00
79) Chrysene-d12	21.269	240	859255	20.000	ng/ul	0.00
88) Perylene-d12	24.168	264	1109463	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.146	96	7863m	3.461	ng/uL	0.00
4) Pyridine-d5	3.564	84	41131	6.038	ng/ul	0.00
7) Phenol-d5	6.816	99	47043	5.459	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.981	67	170299	31.995	ng/ul	0.00
11) 2-Chlorophenol-d4	7.175	132	156891	23.999	ng/ul	0.00
15) 4-Methylphenol-d8	8.346	113	102250	14.157	ng/ul	0.00
21) Nitrobenzene-d5	8.793	128	107373	31.158	ng/ul	0.00
24) 2-Nitrophenol-d4	9.510	143	97507	30.930	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.040	165	195512	28.328	ng/ul	0.00
31) 4-Chloroaniline-d4	10.563	131	267054	23.504	ng/ul	0.00
46) Dimethylphthalate-d6	13.698	166	842003	35.741	ng/ul	0.00
49) Acenaphthylene-d8	13.975	160	912514	33.688	ng/ul	0.00
54) 4-Nitrophenol-d4	14.498	143	16463	3.653	ng/ul	0.00
60) Fluorene-d10	15.281	176	728837	35.354	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.404	200	84059	23.590	ng/ul	0.00
73) Anthracene-d10	17.134	188	1180876	33.950	ng/ul	0.00
81) Pyrene-d10	19.433	212	1557369	33.711	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.968	264	1931508	37.783	ng/ul	0.00

Target Compounds Qvalue

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

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