

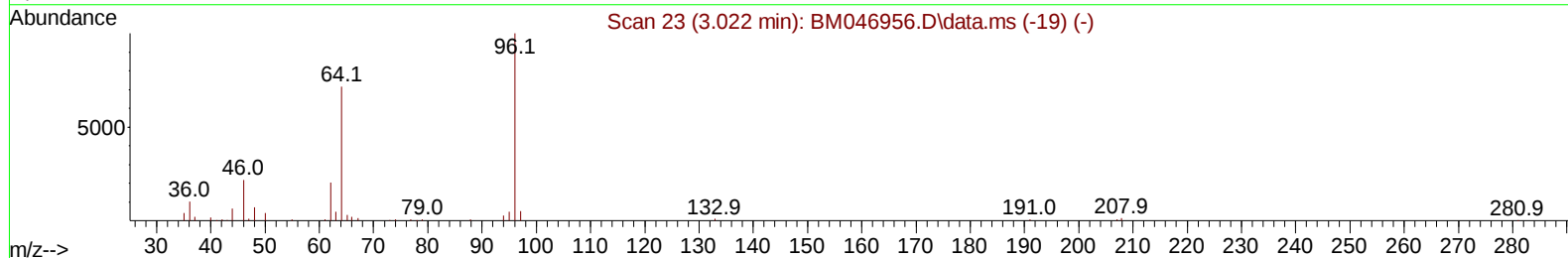
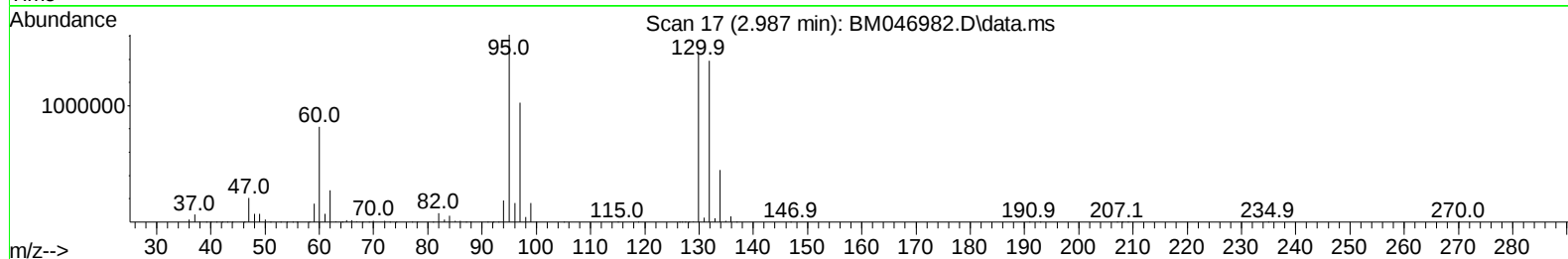
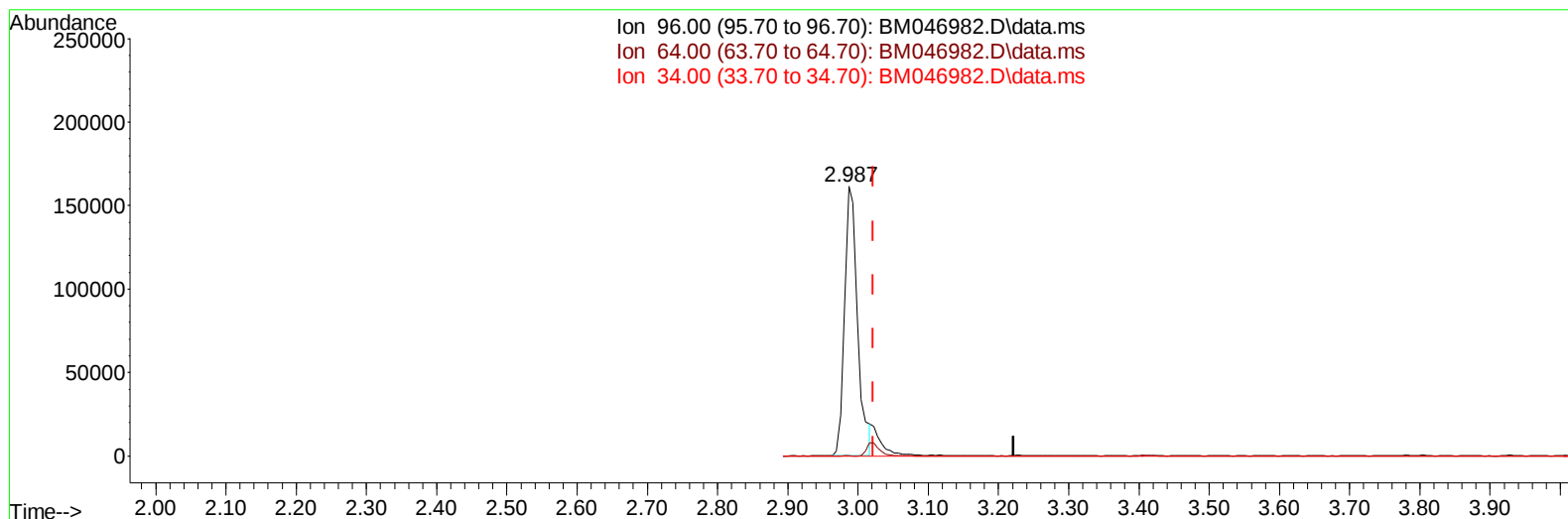
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080324\
Data File : BM046982.D
Acq On : 04 Aug 2024 02:51
Operator : MA/JU
Sample : P3276-23
Misc :
ALS Vial : 29 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E20L1

Manual IntegrationsAPPROVED

Quant Time: Aug 05 00:59:34 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Aug 03 02:51:11 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 08/05/2024
Supervised By :mohammad ahmed 08/05/2024



TIC: BM046982.D\data.ms

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

2.987min (-0.035) 42.58 ng/uL

response 206344

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	81.60	0.18#
34.00	0.00	0.00
0.00	0.00	0.00

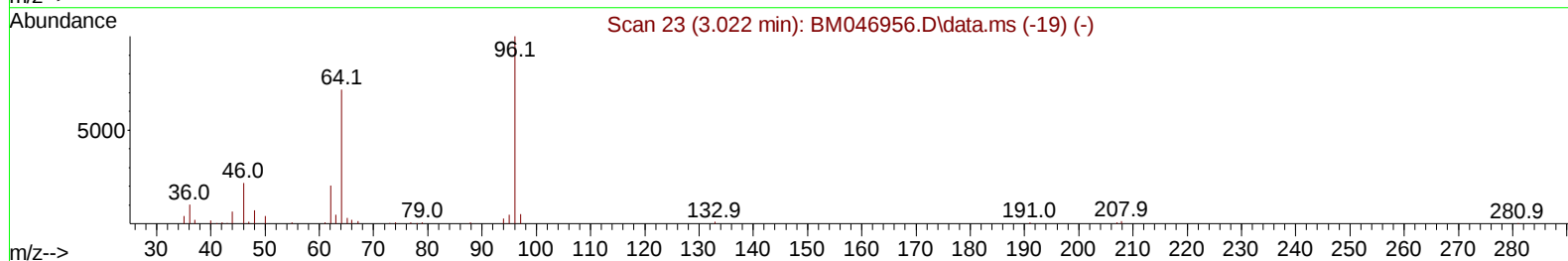
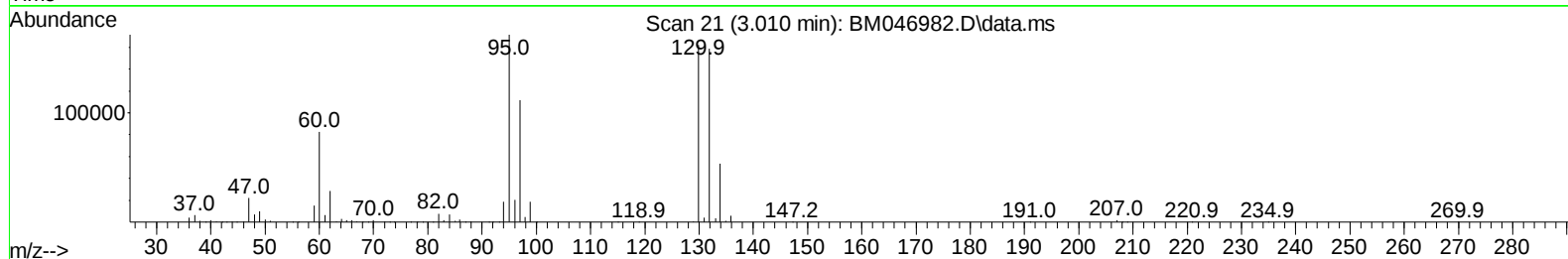
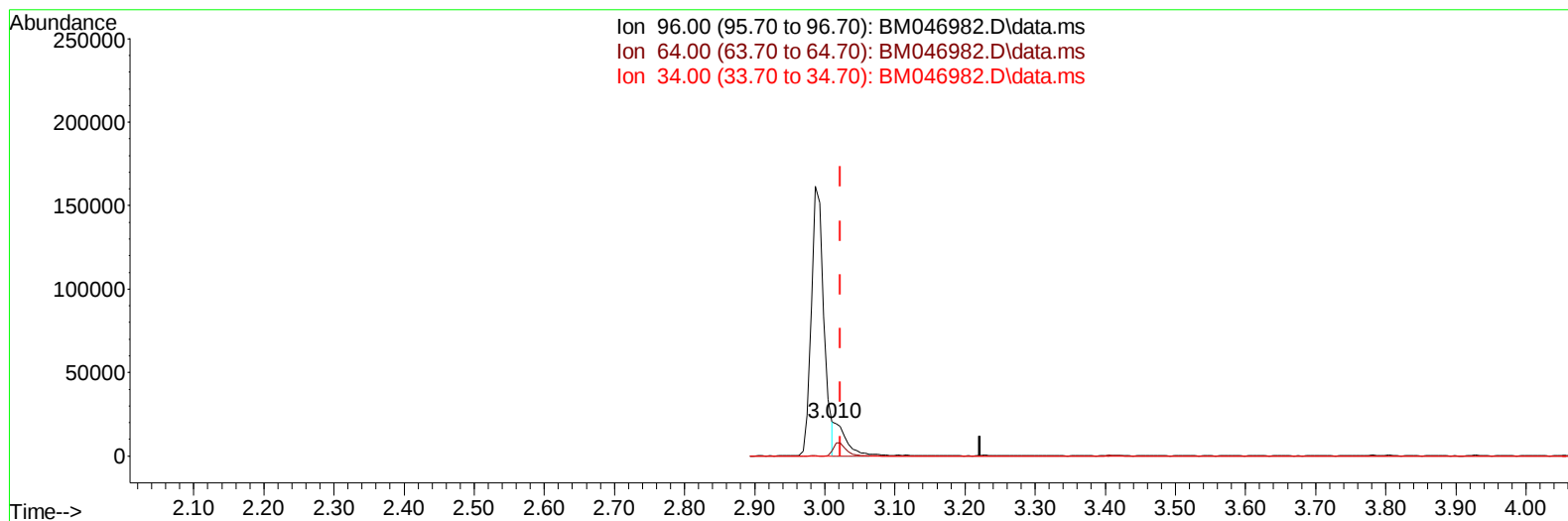
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TIC: BM046982.D\data.ms

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

3.010min (-0.012) 4.95 ng/uL m

response 23963

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	81.60	14.07#
34.00	0.00	0.00
0.00	0.00	0.00

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Quant Time: Aug 05 01:45:14 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.398	152		191973	20.000	ng/ul	0.00
20) Naphthalene-d8	10.157	136		737144	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.057	164		452958	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.815	188		942783	20.000	ng/ul	0.00
79) Chrysene-d12	21.050	240		889504	20.000	ng/ul	0.00
88) Perylene-d12	23.762	264		1090856	20.000	ng/ul	-0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.010	96		23963m	4.945	ng/uL	-0.01
4) Pyridine-d5	3.410	84		224178	16.961	ng/ul	0.00
7) Phenol-d5	6.598	99		275488	18.891	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.751	67		178615	18.615	ng/ul	0.00
11) 2-Chlorophenol-d4	6.940	132		224949	18.210	ng/ul	0.00
15) 4-Methylphenol-d8	8.116	113		214931	18.605	ng/ul	0.00
21) Nitrobenzene-d5	8.539	128		104872	17.729	ng/ul	0.00
24) 2-Nitrophenol-d4	9.257	143		114751	17.764	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.792	165		211232	17.032	ng/ul	0.00
31) 4-Chloroaniline-d4	10.304	131		287816	17.658	ng/ul	0.00
46) Dimethylphthalate-d6	13.480	166		643770	18.588	ng/ul	0.00
49) Acenaphthylene-d8	13.739	160		721000	18.650	ng/ul	0.00
54) 4-Nitrophenol-d4	14.292	143		88464	13.211	ng/ul	0.00
60) Fluorene-d10	15.057	176		532983	17.990	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.198	200		70786	11.884	ng/ul	0.00
73) Anthracene-d10	16.909	188		840405	18.804	ng/ul	0.00
81) Pyrene-d10	19.227	212		961549	18.884	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.574	264		1081945	18.836	ng/ul	-0.01

Target Compounds Qvalue

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

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