

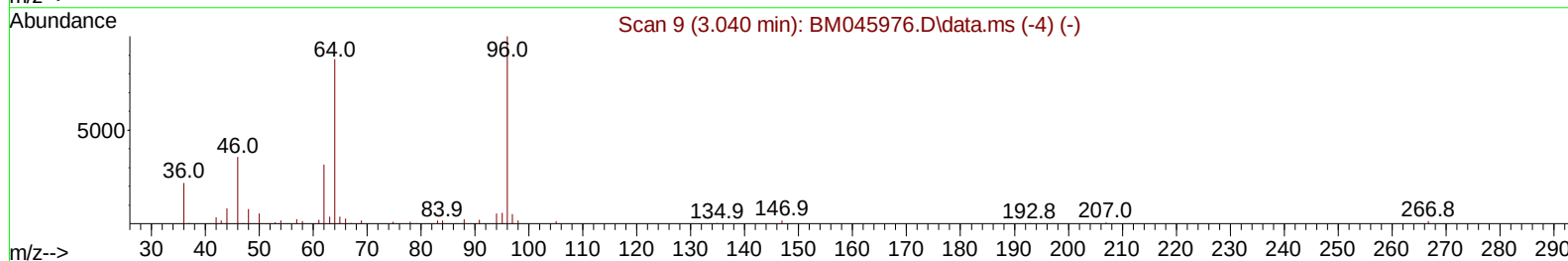
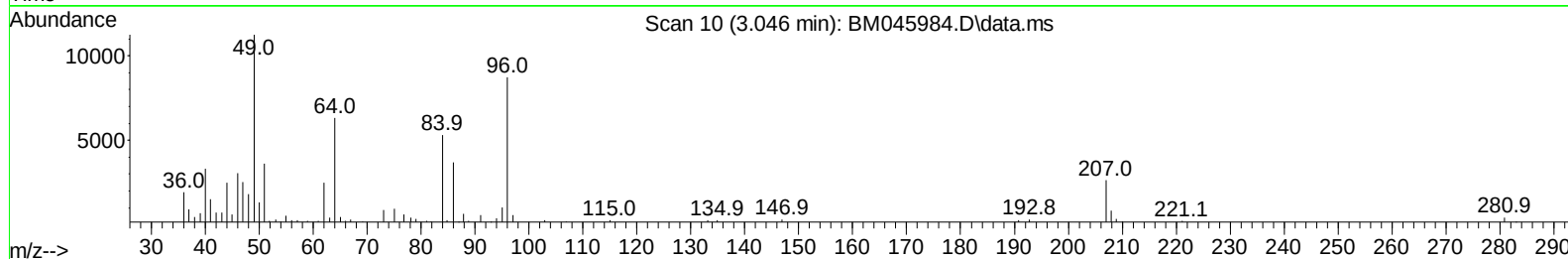
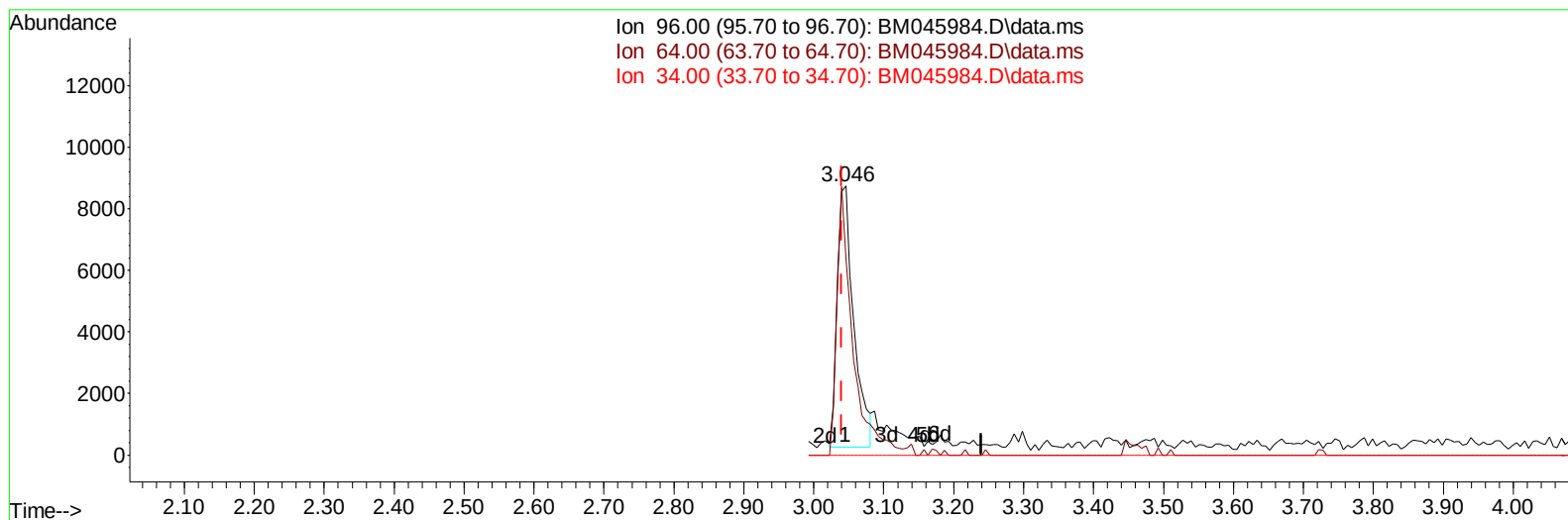
Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM061124\
Data File : BM045984.D
Acq On : 11 Jun 2024 20:19
Operator : MA/JU
Sample : P2642-07
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
A4BH2

Manual IntegrationsAPPROVED

Quant Time: Jun 11 23:06:40 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 11 16:01:34 2024
Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 06/12/2024
Supervised By :mohammad ahmed 06/13/2024



TIC: BM045984.D\data.ms

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

3.046min (+ 0.006) 4.01 ng/uL

response 14127

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	88.70	72.64
34.00	0.00	0.00
0.00	0.00	0.00

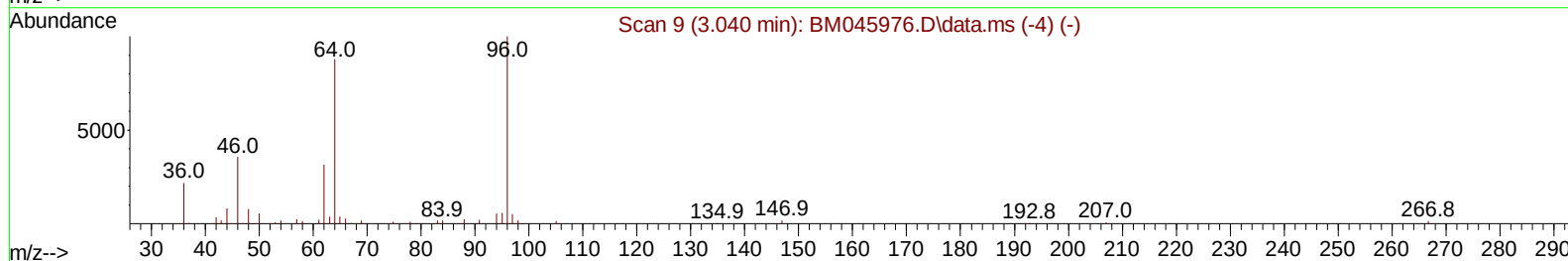
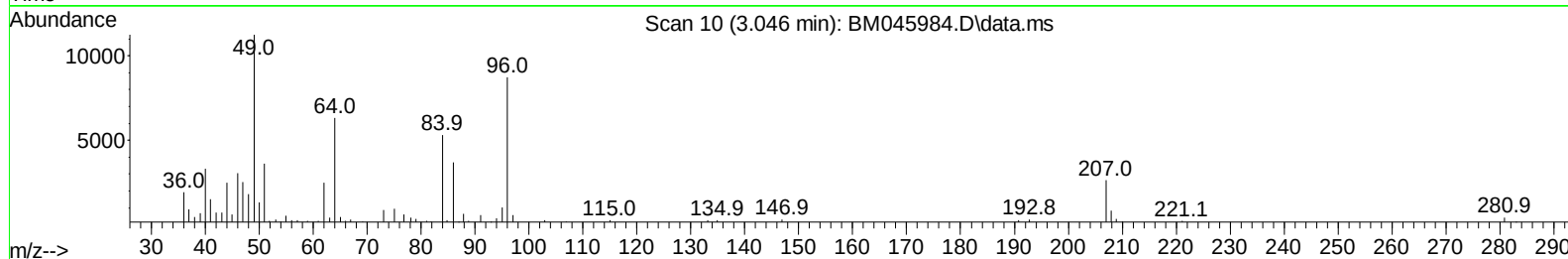
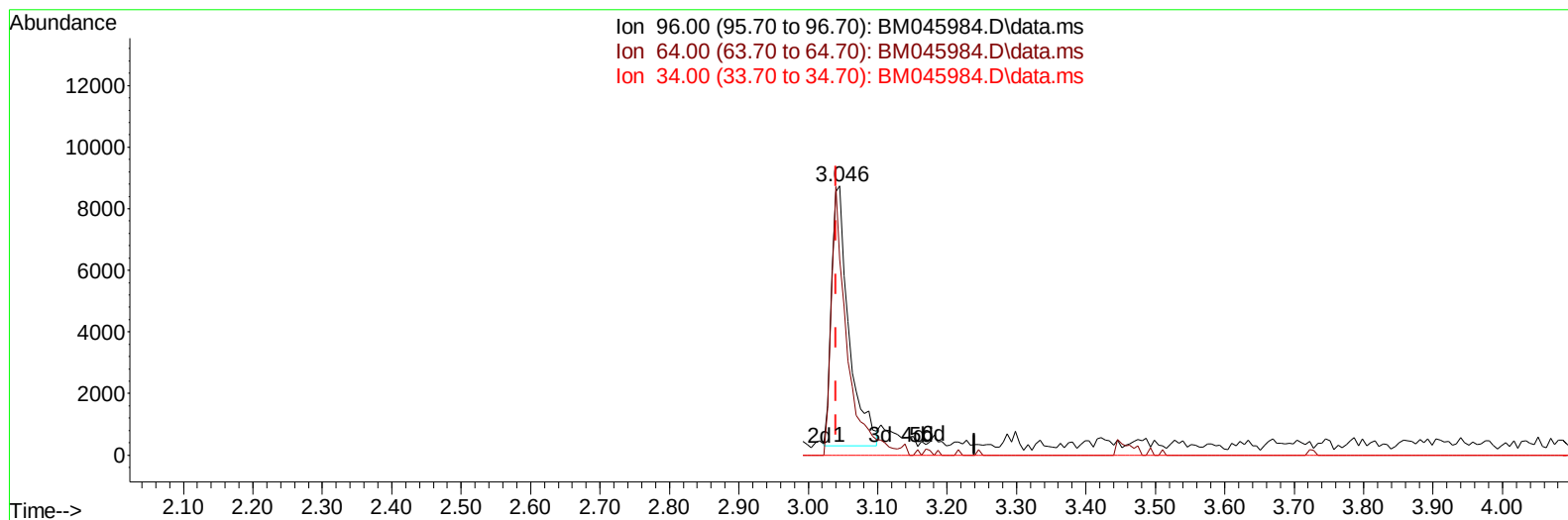
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TIC: BM045984.D\data.ms

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

3.046min (+ 0.006) 4.17 ng/uL m

response 14708

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	88.70	72.64
34.00	0.00	0.00
0.00	0.00	0.00

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 ALS Vial : 10 Sample Multiplier: 1

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Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 06/12/2024
 Supervised By :mohammad ahmed 06/13/2024

Quant Time: Jun 11 23:18:02 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM053124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Jun 11 16:01:34 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.434	152	136719	20.000	ng/ul	0.00
20) Naphthalene-d8	10.204	136	603134	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.086	164	397268	20.000	ng/ul	0.00
64) Phenanthrene-d10	16.827	188	855404	20.000	ng/ul	0.00
79) Chrysene-d12	21.044	240	879004	20.000	ng/ul	0.00
88) Perylene-d12	23.744	264	929716	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.046	96	14708m	4.171	ng/uL	0.00
4) Pyridine-d5	3.452	84	196188	21.607	ng/ul	0.00
7) Phenol-d5	6.692	99	310534	26.736	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	6.804	67	225401	27.026	ng/ul	0.00
11) 2-Chlorophenol-d4	6.998	132	242427	27.943	ng/ul	0.00
15) 4-Methylphenol-d8	8.204	113	255268	29.091	ng/ul	0.00
21) Nitrobenzene-d5	8.604	128	125803	26.678	ng/ul	0.00
24) 2-Nitrophenol-d4	9.316	143	125530	26.302	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	9.863	165	235504	26.314	ng/ul	0.00
31) 4-Chloroaniline-d4	10.386	131	339960	27.446	ng/ul	0.00
46) Dimethylphthalate-d6	13.533	166	793704	28.859	ng/ul	0.00
49) Acenaphthylene-d8	13.774	160	904205	28.312	ng/ul	0.00
54) 4-Nitrophenol-d4	14.386	143	125699	21.177	ng/ul	0.00
60) Fluorene-d10	15.086	176	686929	29.277	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.227	200	84105	19.254	ng/ul	0.00
73) Anthracene-d10	16.927	188	1081213	29.075	ng/ul	0.00
81) Pyrene-d10	19.227	212	1360196	24.883	ng/ul	0.00
92) Benzo(a)pyrene-d12	23.562	264	1316913	30.307	ng/ul	0.00

Target Compounds Qvalue

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

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