

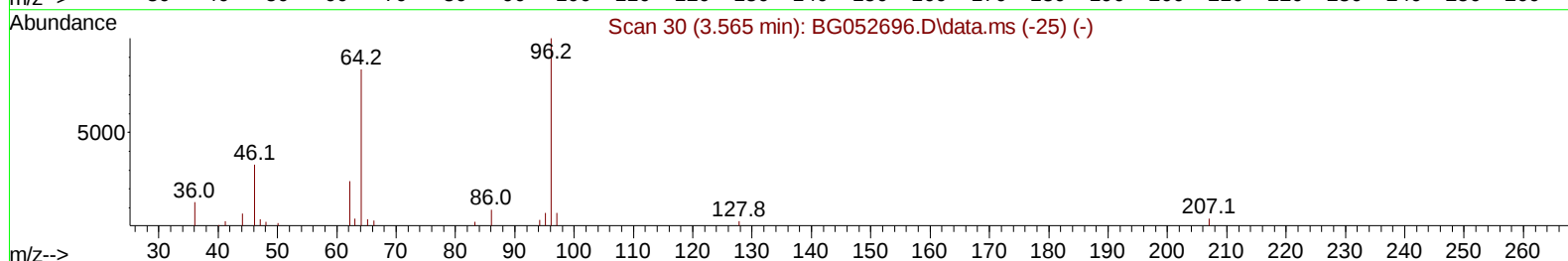
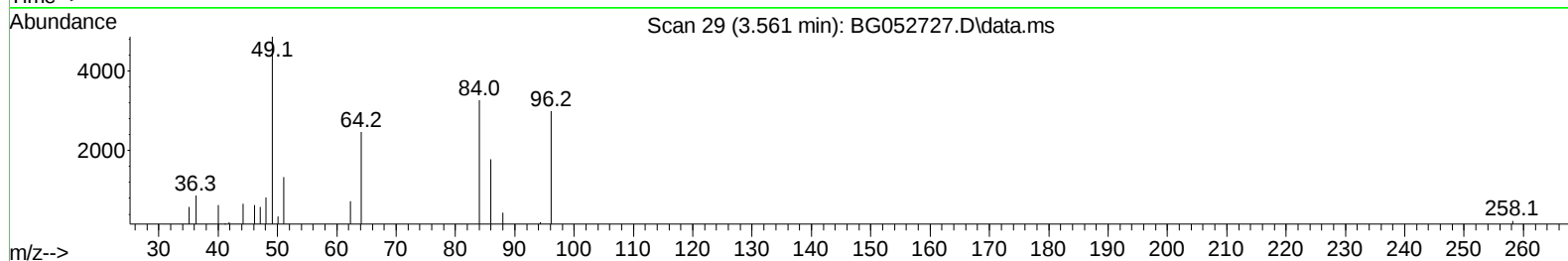
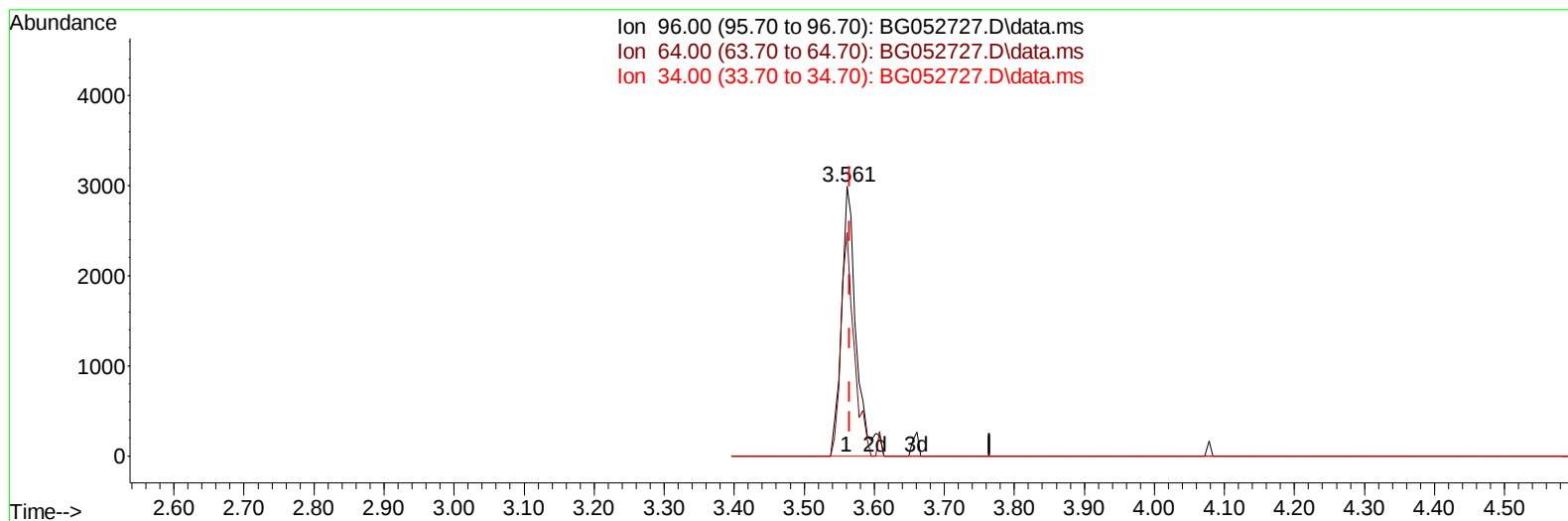
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031822\
Data File : BG052727.D
Acq On : 19 Mar 2022 2:09
Operator : CG/JU
Sample : N1882-02
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
DBRX4

Manual IntegrationsAPPROVED

Reviewed By :Jagrut Upadhyay 03/21/2022
Supervised By :Yogesh Patel 03/24/2022

Quant Time: Mar 19 03:29:30 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Mar 18 00:55:42 2022
Response via : Initial Calibration



TIC: BG052727.D\data.ms

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

3.561min (-0.004) 4.11 ng/uL

response 4244

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	82.81
34.00	0.00	0.00
0.00	0.00	0.00

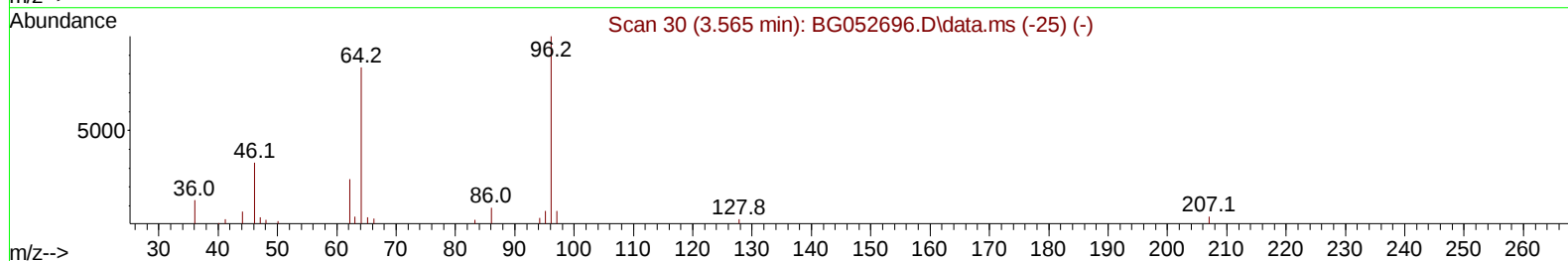
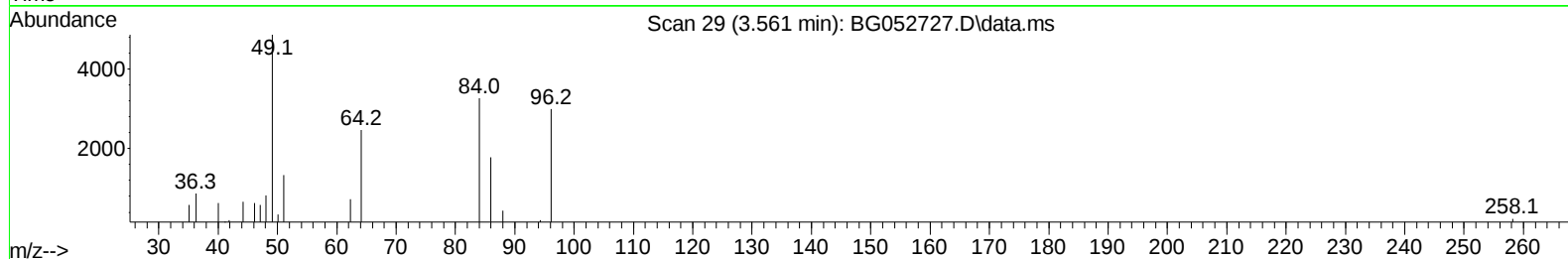
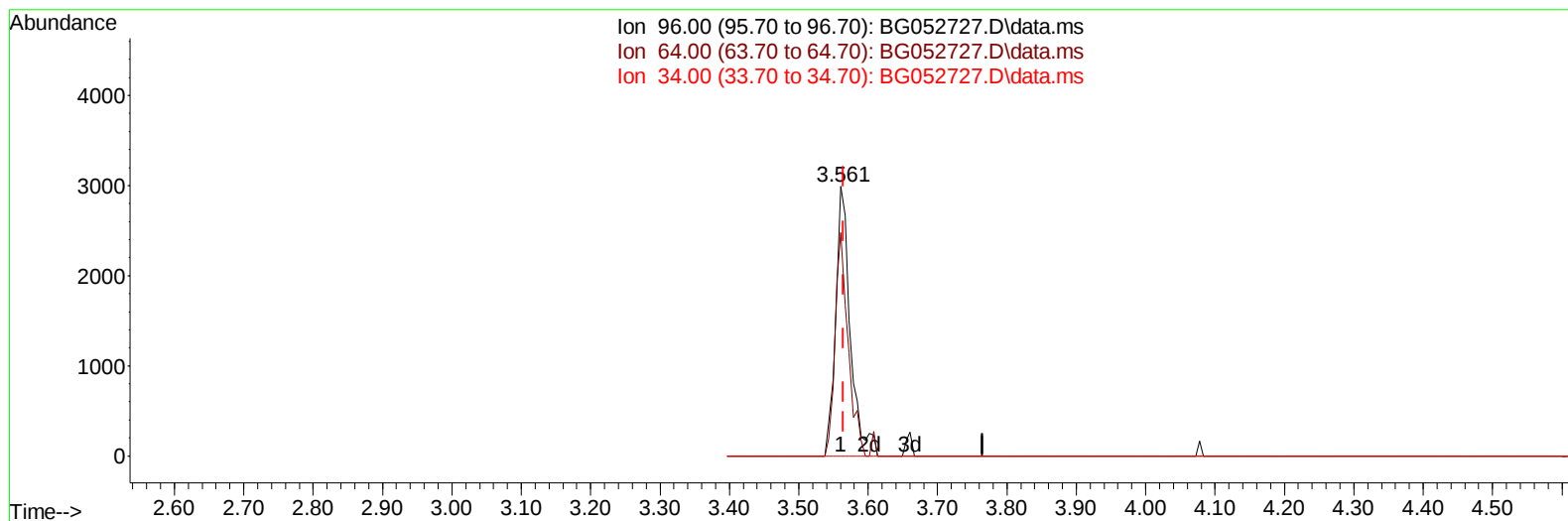
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TIC: BG052727.D\data.ms

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

3.561min (-0.004) 4.28 ng/uL m

response 4413

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	82.81
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	8.149	152	36457	20.000	ng/ul	0.00
20) Naphthalene-d8	10.963	136	149576	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.769	164	105720	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.512	188	247974	20.000	ng/ul	0.00
79) Chrysene-d12	21.795	240	276954	20.000	ng/ul	-0.01
88) Perylene-d12	25.114	264	273302	20.000	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.561	96	4413m	4.276	ng/uL	0.00
4) Pyridine-d5	3.978	84	56367	21.876	ng/ul	0.00
7) Phenol-d5	7.291	99	75305	23.062	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.467	67	46884	23.509	ng/ul	0.00
11) 2-Chlorophenol-d4	7.679	132	52625	23.735	ng/ul	0.00
15) 4-Methylphenol-d8	8.836	113	54348	22.977	ng/ul	-0.01
21) Nitrobenzene-d5	9.306	128	24844	23.074	ng/ul	0.00
24) 2-Nitrophenol-d4	10.034	143	27029	23.800	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.569	165	52072	21.077	ng/ul	0.00
31) 4-Chloroaniline-d4	11.086	131	72429	22.181	ng/ul	0.00
46) Dimethylphthalate-d6	14.164	166	184400	22.704	ng/ul	0.00
49) Acenaphthylene-d8	14.458	160	227432	23.044	ng/ul	-0.01
54) 4-Nitrophenol-d4	14.928	143	26388	19.324	ng/ul	0.00
60) Fluorene-d10	15.756	176	163206	22.886	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.856	200	22929	18.170	ng/ul	0.00
73) Anthracene-d10	17.606	188	269832	23.368	ng/ul	0.00
81) Pyrene-d10	19.892	212	336105	21.693	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.885	264	325626	23.059	ng/ul	-0.01

Target Compounds Qvalue

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

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