

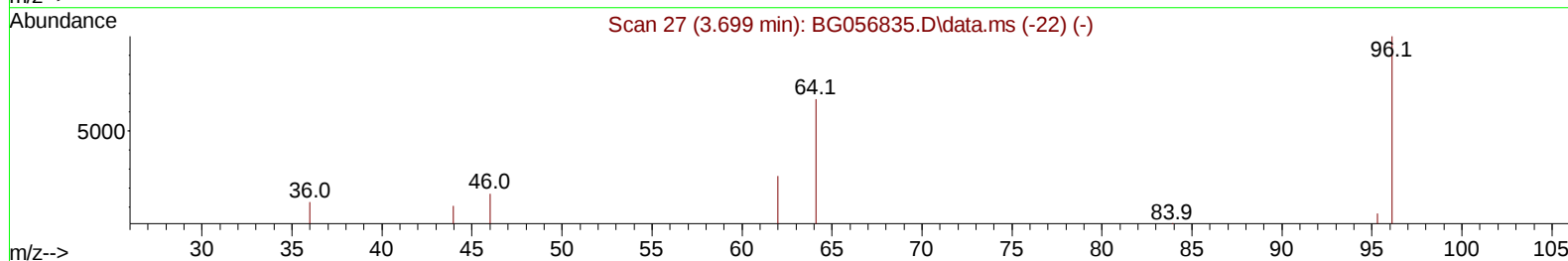
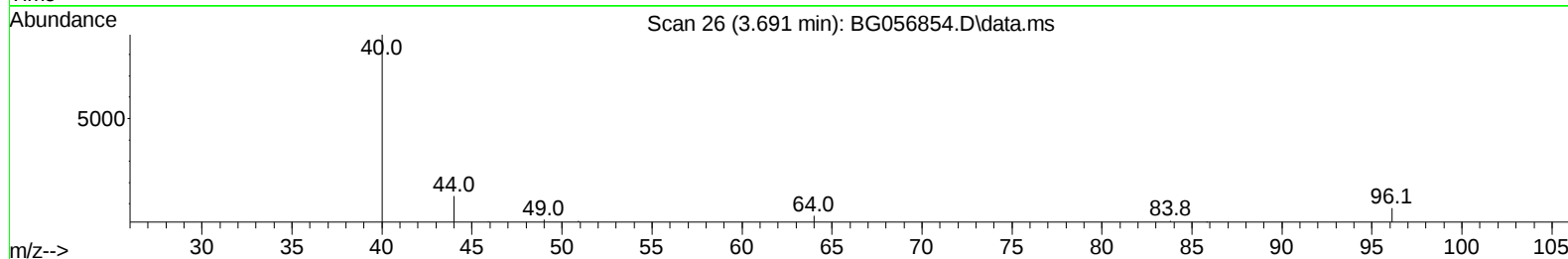
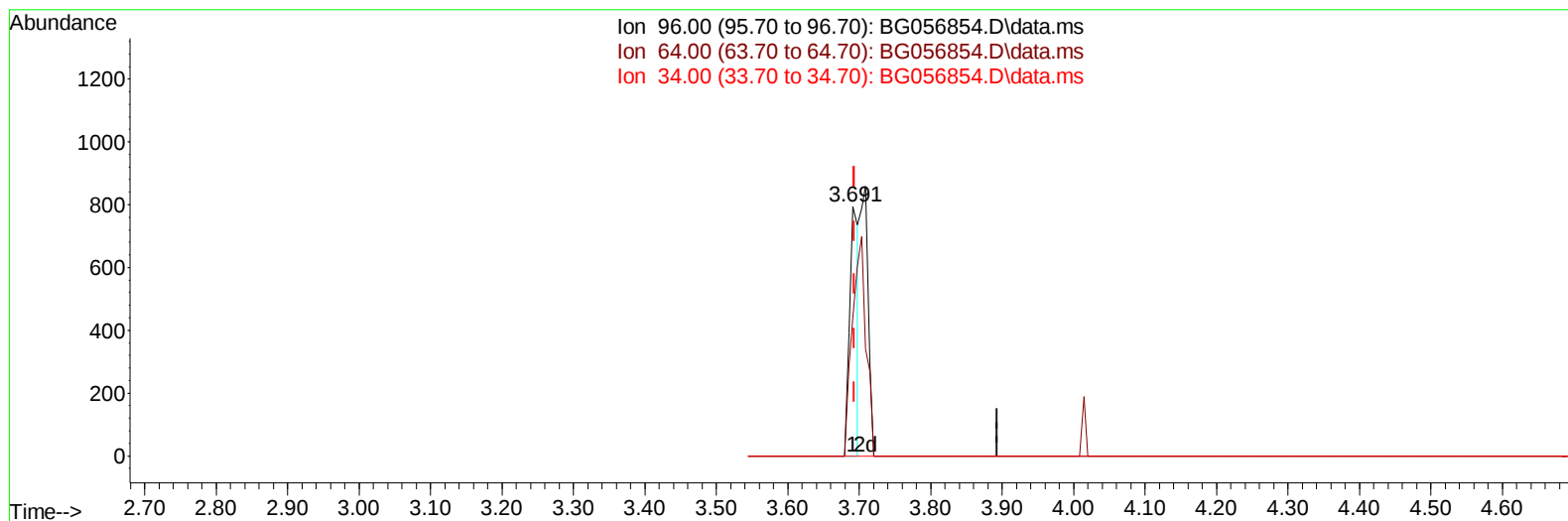
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG030623\
 Data File : BG056854.D
 Acq On : 6 Mar 2023 23:37
 Operator : CG/JU
 Sample : 01712-01
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBAD6

Manual IntegrationsAPPROVED

Reviewed By :Christian Giraldo 03/07/2023
 Supervised By :Jagrut Upadhyay 03/07/2023

Quant Time: Mar 07 00:55:45 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG022323.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Mar 02 00:12:28 2023
 Response via : Initial Calibration



TIC: BG056854.D\data.ms

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

3.691min (-0.002) 2.13 ng/uL

response 678

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	60.40	55.29
34.00	0.00	0.00
0.00	0.00	0.00

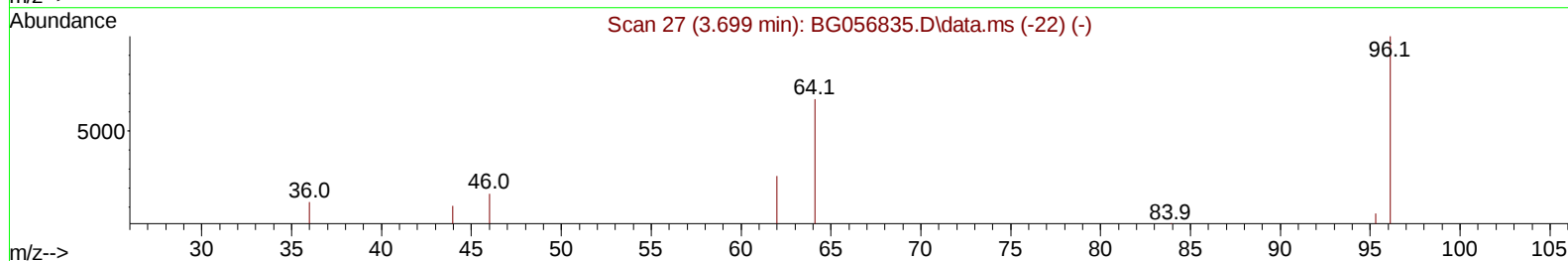
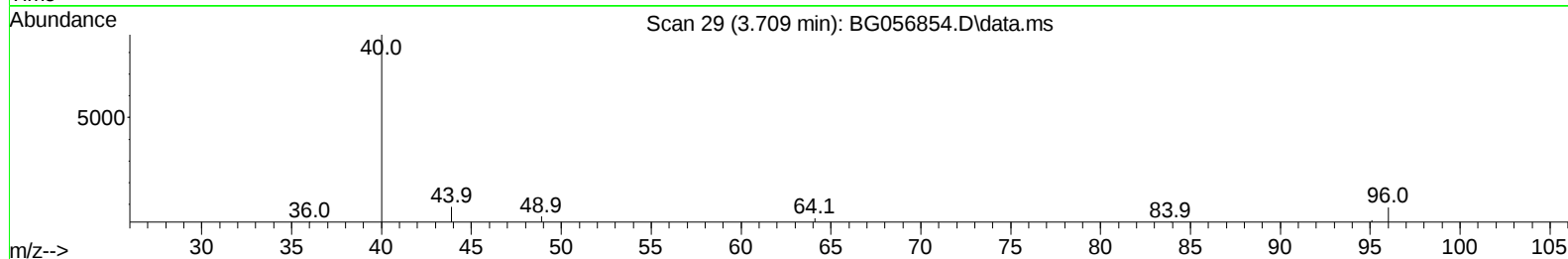
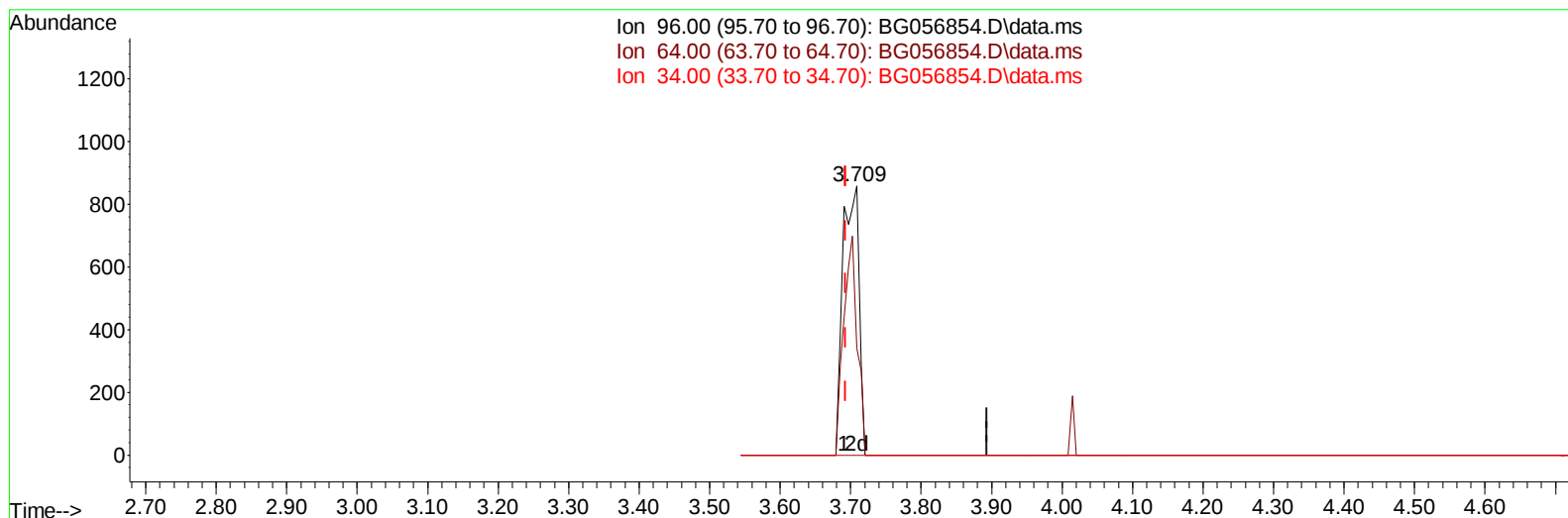
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(3) 1,4-Dioxane-d8 (S)

3.709min (+ 0.015) 4.26 ng/uL m

response 1358

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	60.40	39.63#
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	8.421	152	17670	20.000	ng/ul	0.00
20) Naphthalene-d8	11.253	136	79598	20.000	ng/ul	0.00
38) Acenaphthene-d10	15.019	164	59465	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.751	188	135033	20.000	ng/ul	0.00
79) Chrysene-d12	22.064	240	122132	20.000	ng/ul	0.00
88) Perylene-d12	25.589	264	129221	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.709	96	1358m	4.263	ng/uL	0.02
4) Pyridine-d5	4.138	84	15890	12.359	ng/ul	0.00
7) Phenol-d5	7.540	99	29589	16.734	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.722	67	16154	14.977	ng/ul	0.00
11) 2-Chlorophenol-d4	7.939	132	18971	16.156	ng/ul	0.00
15) 4-Methylphenol-d8	9.102	113	22054	16.323	ng/ul	0.00
21) Nitrobenzene-d5	9.590	128	9861	14.783	ng/ul	0.00
24) 2-Nitrophenol-d4	10.319	143	11535	14.527	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.859	165	22235	14.864	ng/ul	0.00
31) 4-Chloroaniline-d4	11.376	131	22190	11.373	ng/ul	0.00
46) Dimethylphthalate-d6	14.408	166	68979	14.377	ng/ul	0.00
49) Acenaphthylene-d8	14.719	160	79391	14.559	ng/ul	0.00
54) 4-Nitrophenol-d4	15.178	143	8444	10.208	ng/ul	0.00
60) Fluorene-d10	16.000	176	62367	14.053	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	16.106	200	7558	8.366	ng/ul	0.00
73) Anthracene-d10	17.851	188	97261	14.847	ng/ul	0.00
81) Pyrene-d10	20.113	212	115906	14.025	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.342	264	107015	15.343	ng/ul	0.00

Target Compounds Qvalue

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

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