

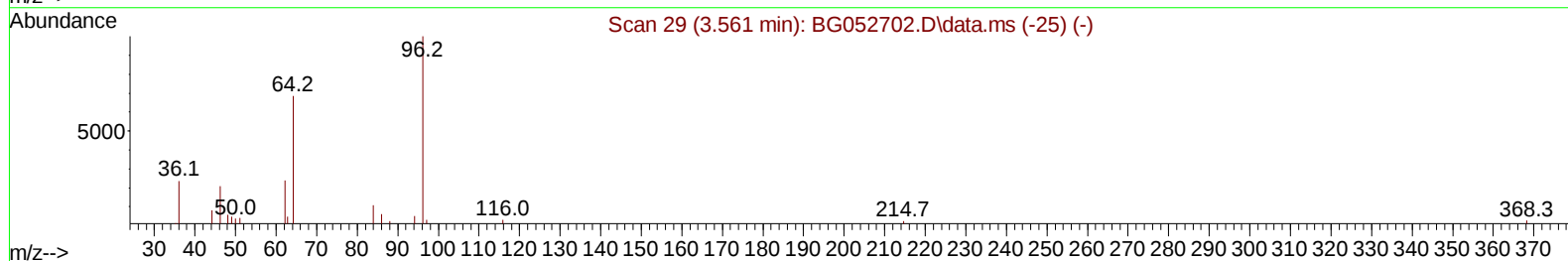
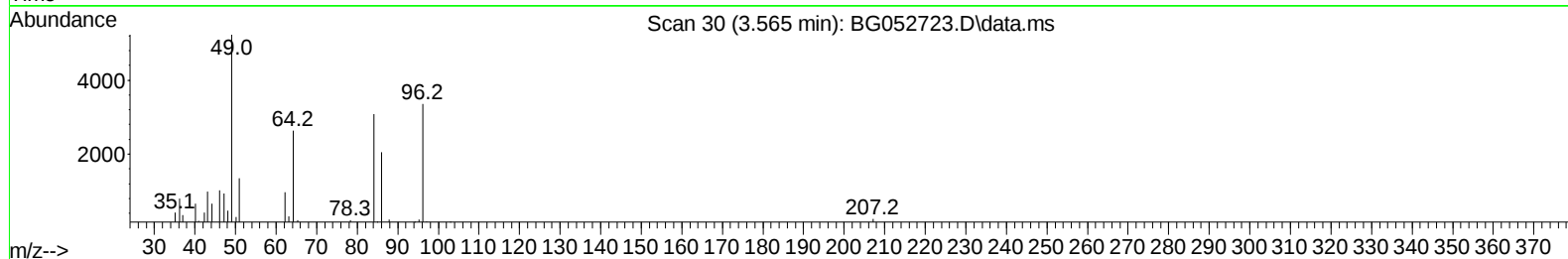
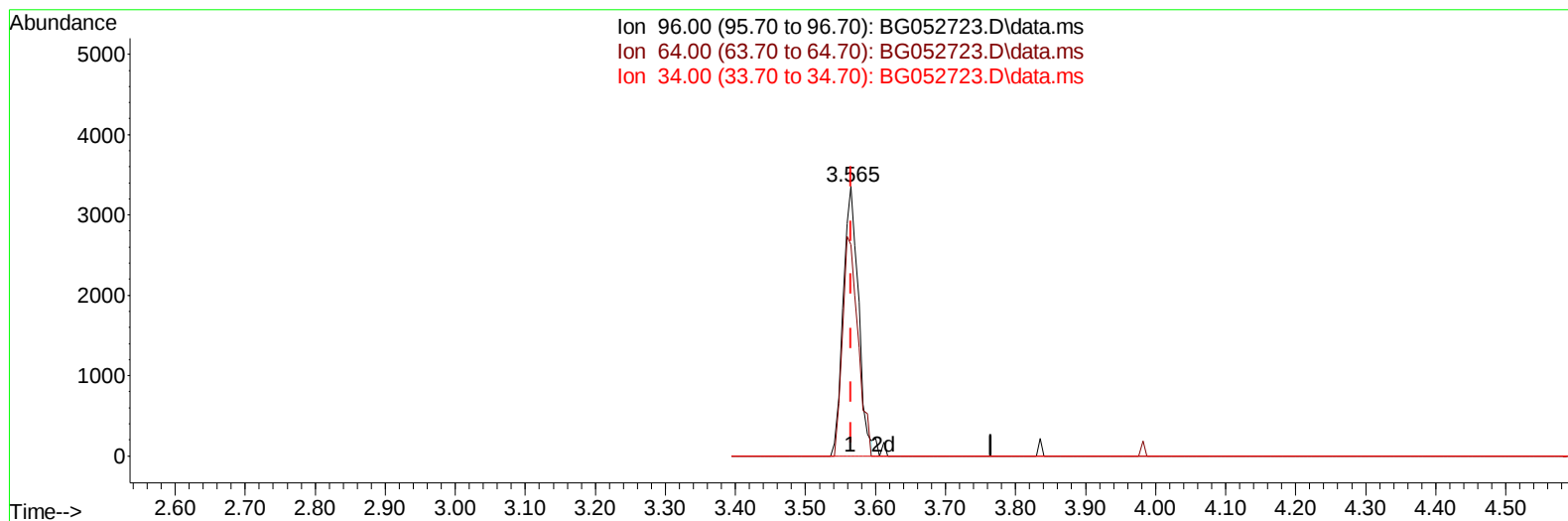
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG031822\
 Data File : BG052723.D
 Acq On : 18 Mar 2022 23:33
 Operator : CG/JU
 Sample : N1882-07
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 DBRX7

Manual IntegrationsAPPROVED

Quant Time: Mar 19 01:01:17 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

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



TIC: BG052723.D\data.ms

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

3.565min (-0.000) 5.77 ng/uL

response 5152

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	78.36
34.00	0.00	0.00
0.00	0.00	0.00

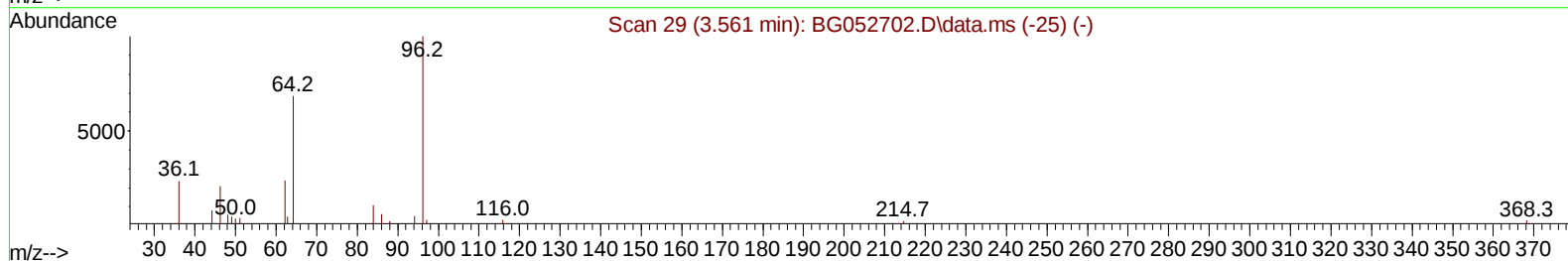
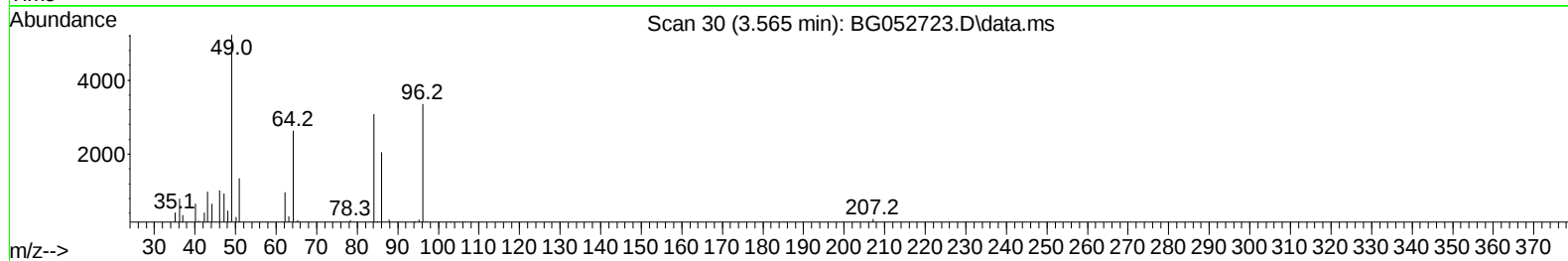
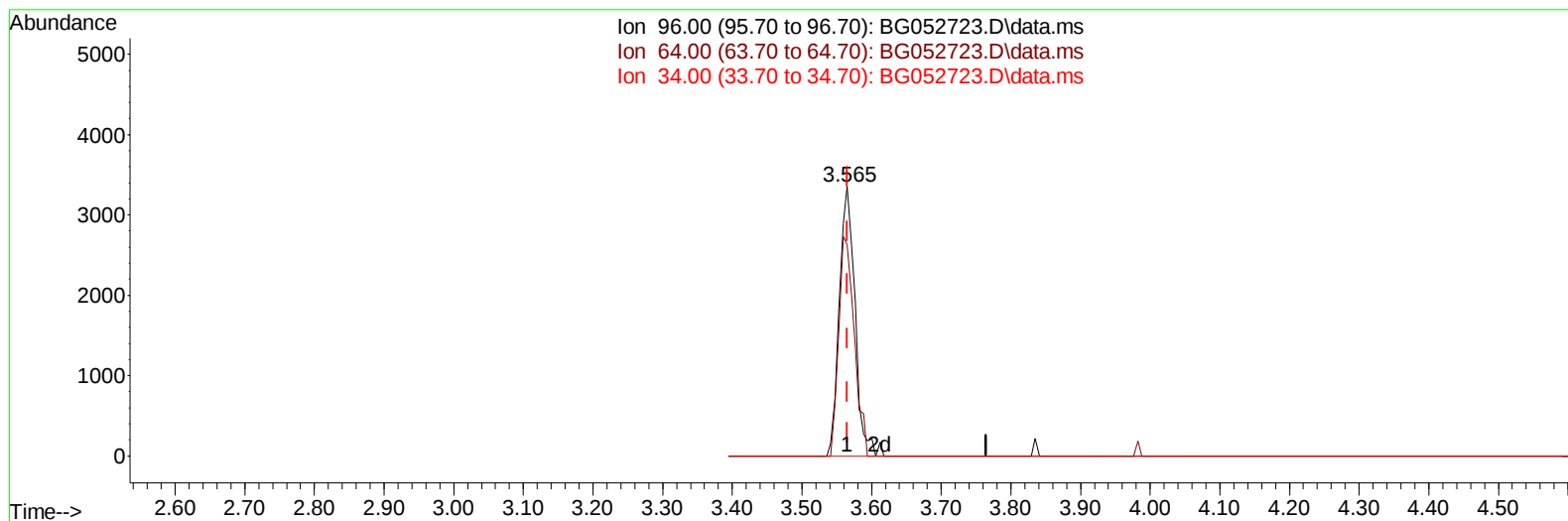
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TIC: BG052723.D\data.ms

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

3.565min (-0.000) 5.86 ng/uL m

response 5230

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	78.36
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	8.147	152	31521	20.000	ng/ul	0.00
20) Naphthalene-d8	10.961	136	129122	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.767	164	92283	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.511	188	215964	20.000	ng/ul	0.00
79) Chrysene-d12	21.799	240	232405	20.000	ng/ul	0.00
88) Perylene-d12	25.118	264	227585	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.565	96	5230m	5.861	ng/uL	0.00
4) Pyridine-d5	3.982	84	65805	29.538	ng/ul	0.00
7) Phenol-d5	7.289	99	94832	33.590	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.466	67	56679	32.871	ng/ul	0.00
11) 2-Chlorophenol-d4	7.677	132	64022	33.397	ng/ul	0.00
15) 4-Methylphenol-d8	8.840	113	69640	34.052	ng/ul	0.00
21) Nitrobenzene-d5	9.304	128	32075	34.509	ng/ul	0.00
24) 2-Nitrophenol-d4	10.033	143	34290	34.977	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.573	165	66272	31.073	ng/ul	0.00
31) 4-Chloroaniline-d4	11.084	131	87960	31.205	ng/ul	0.00
46) Dimethylphthalate-d6	14.168	166	221443	31.235	ng/ul	0.00
49) Acenaphthylene-d8	14.462	160	271359	31.499	ng/ul	0.00
54) 4-Nitrophenol-d4	14.932	143	35019	29.379	ng/ul	0.00
60) Fluorene-d10	15.754	176	191971	30.839	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.860	200	31005	28.211	ng/ul	0.00
73) Anthracene-d10	17.611	188	317109	31.532	ng/ul	0.00
81) Pyrene-d10	19.890	212	401438	30.877	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.889	264	374974	31.887	ng/ul	0.00

Target Compounds Qvalue

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

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