

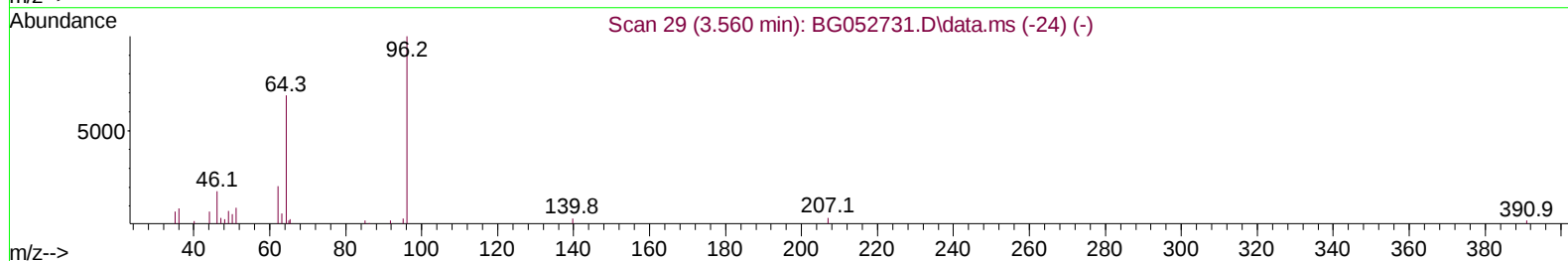
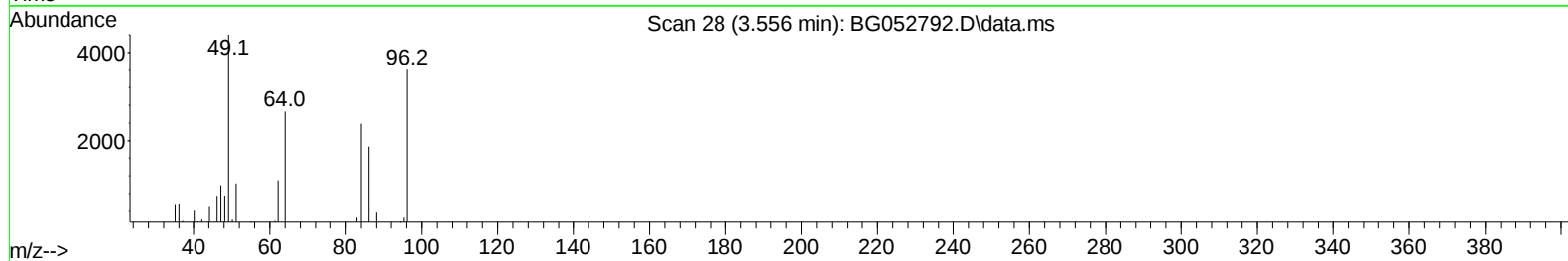
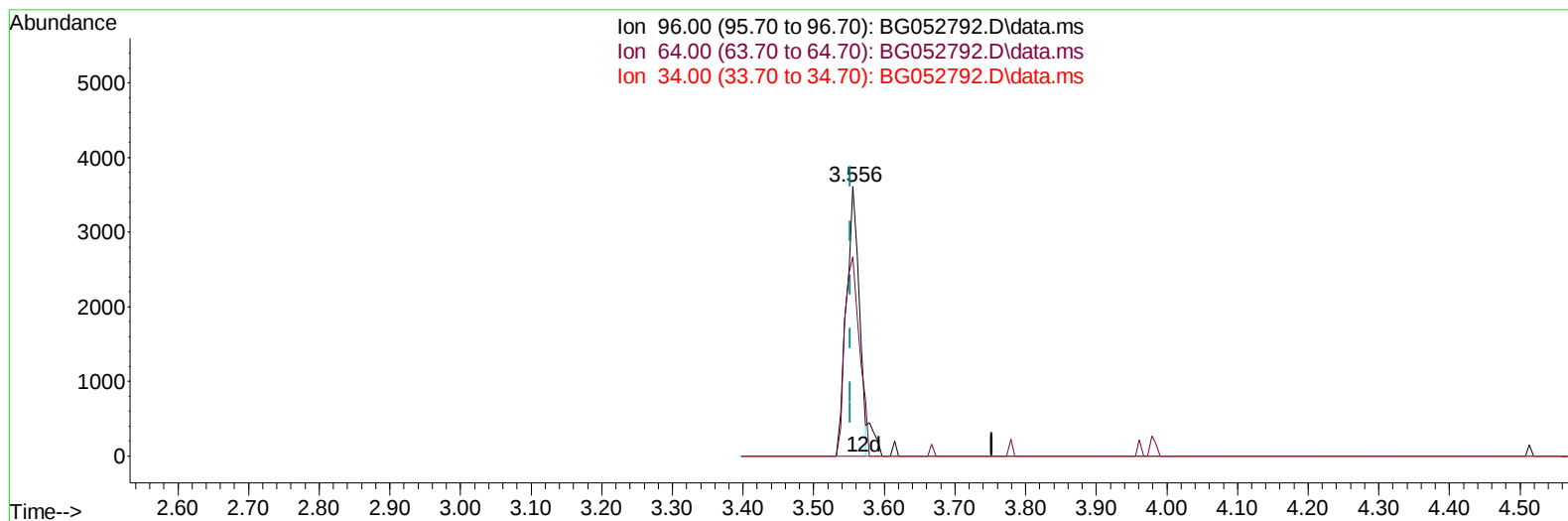
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032322\
 Data File : BG052792.D
 Acq On : 23 Mar 2022 11:16
 Operator : CG/JU
 Sample : PB143280BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK280

Manual IntegrationsAPPROVED

Quant Time: Mar 23 13:24:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG031822.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Mar 23 13:16:32 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 03/23/2022
 Supervised By :mohammad ahmed 03/25/2022



TIC: BG052792.D\data.ms

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

3.556min (+ 0.003) 6.12 ng/uL

response 4614

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	73.82
34.00	0.00	0.00
0.00	0.00	0.00

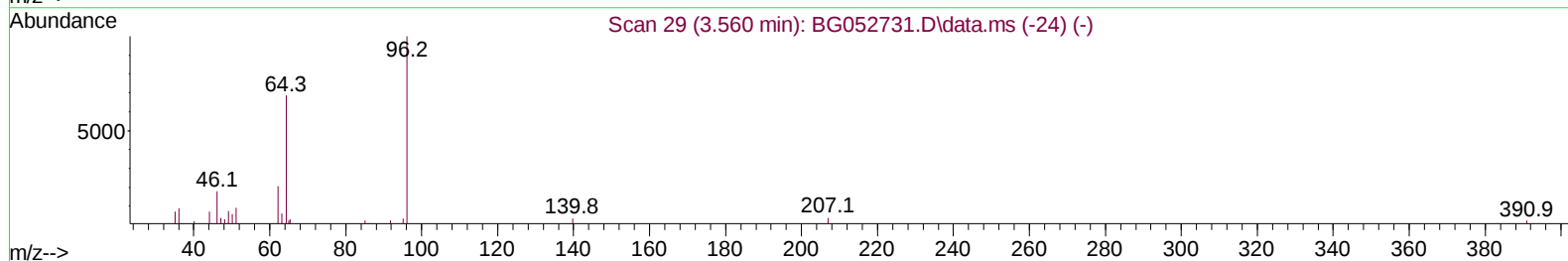
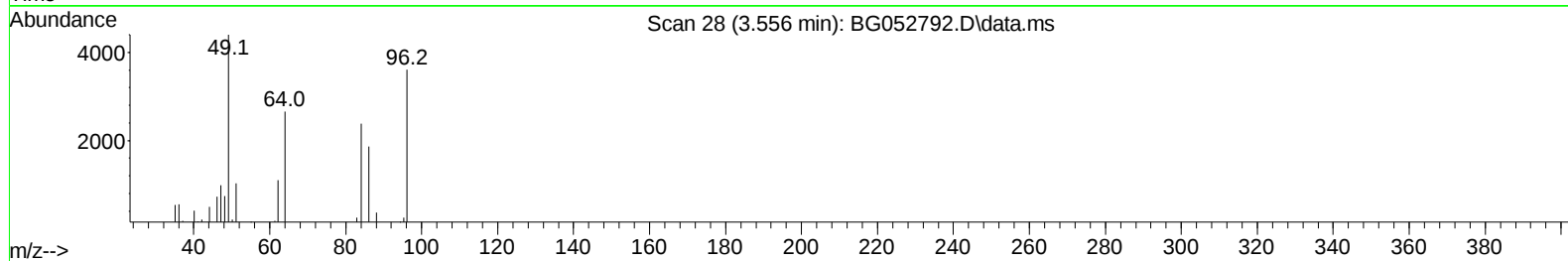
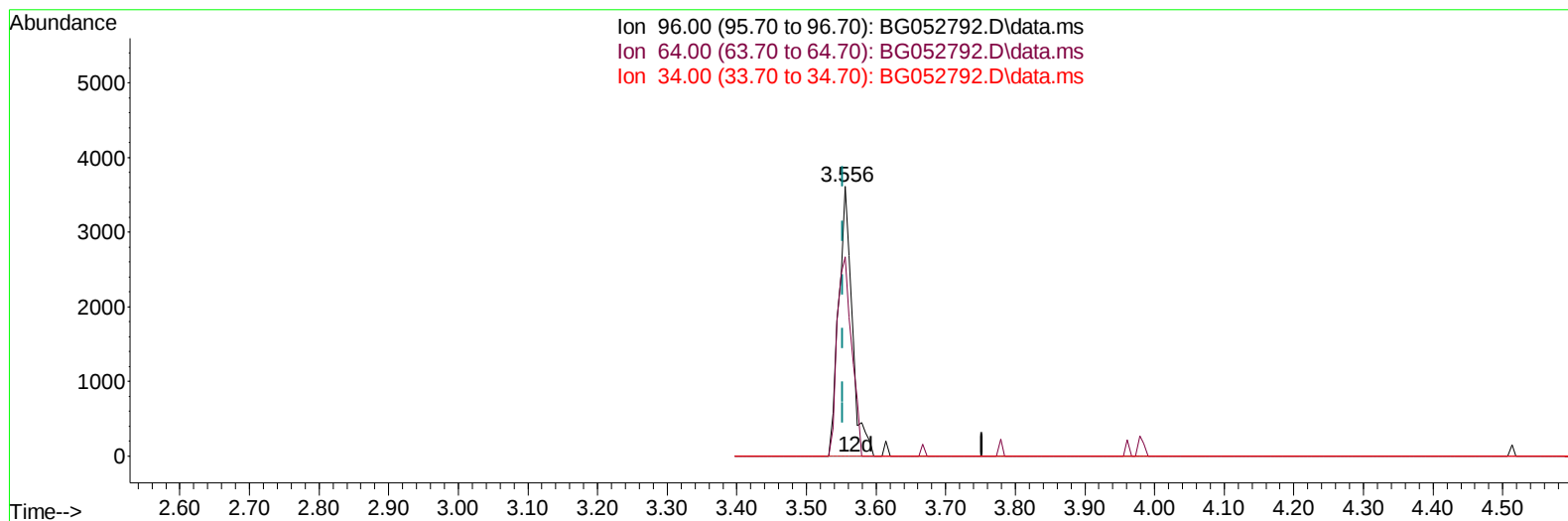
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TIC: BG052792.D\data.ms

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

3.556min (+ 0.003) 6.57 ng/uL m

response 4950

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	73.82
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	8.144	152	26628	20.000	ng/ul	0.00
20) Naphthalene-d8	10.957	136	110677	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.764	164	84835	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.507	188	209327	20.000	ng/ul	0.00
79) Chrysene-d12	21.790	240	211815	20.000	ng/ul	0.00
88) Perylene-d12	25.109	264	201948	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.556	96	4950m	6.566	ng/uL	0.00
4) Pyridine-d5	3.973	84	63510	33.746	ng/ul	0.00
7) Phenol-d5	7.286	99	87355	36.628	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.456	67	54212	37.218	ng/ul	0.00
11) 2-Chlorophenol-d4	7.674	132	59525	36.757	ng/ul	0.00
15) 4-Methylphenol-d8	8.837	113	63931	37.005	ng/ul	0.00
21) Nitrobenzene-d5	9.301	128	31324	39.317	ng/ul	0.00
24) 2-Nitrophenol-d4	10.023	143	35786	42.586	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.570	165	64255	35.148	ng/ul	0.00
31) 4-Chloroaniline-d4	11.081	131	91417	37.836	ng/ul	0.00
46) Dimethylphthalate-d6	14.159	166	260576	39.982	ng/ul	0.00
49) Acenaphthylene-d8	14.459	160	288760	36.461	ng/ul	0.00
54) 4-Nitrophenol-d4	14.928	143	43078	39.313	ng/ul	0.00
60) Fluorene-d10	15.751	176	214940	37.560	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.857	200	40150	37.690	ng/ul	0.00
73) Anthracene-d10	17.601	188	367480	37.700	ng/ul	0.00
81) Pyrene-d10	19.886	212	448546	37.854	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.886	264	396959	38.042	ng/ul	0.00

Target Compounds Qvalue

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

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