

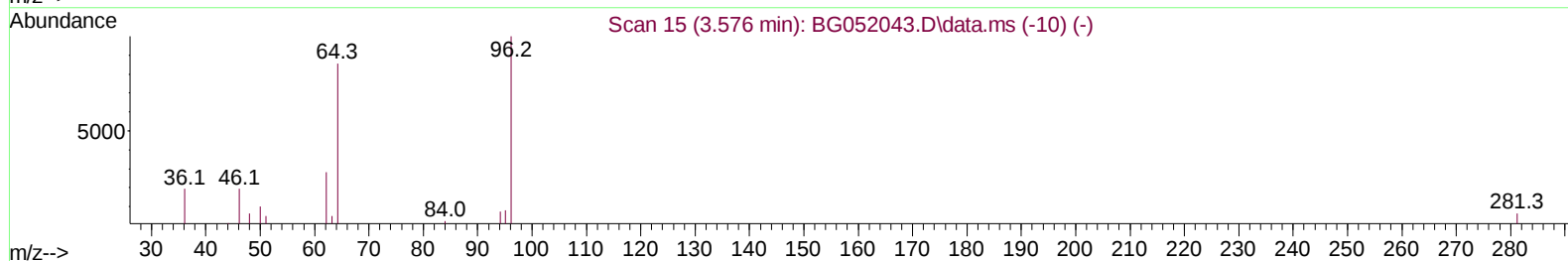
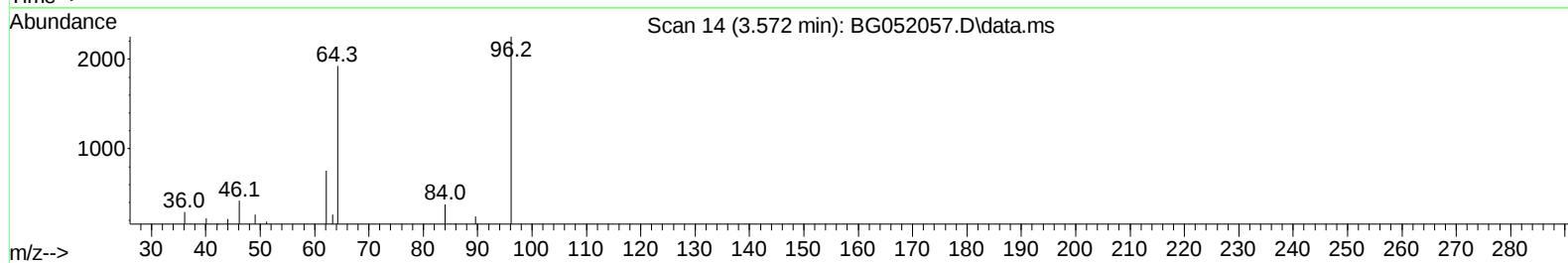
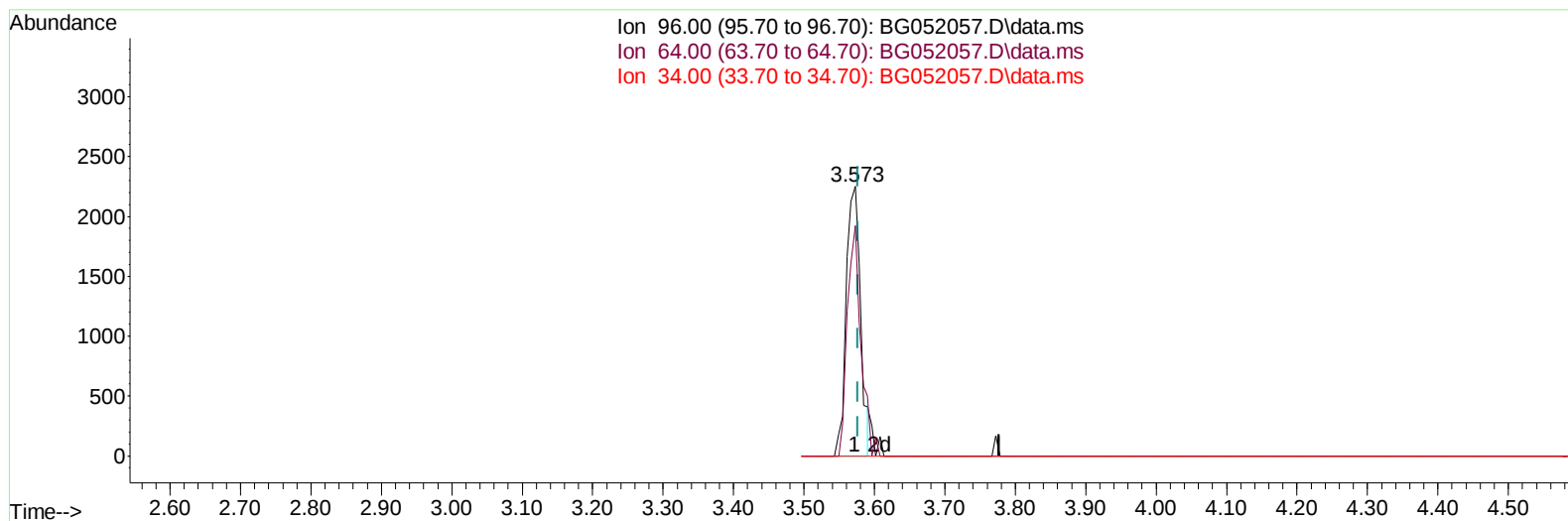
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011722\
 Data File : BG052057.D
 Acq On : 18 Jan 2022 18:05
 Operator : CG/JU
 Sample : PB142117BL
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK117

Manual IntegrationsAPPROVED

Quant Time: Jan 19 00:31:32 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 06 14:43:02 2022
 Response via : Initial Calibration

Reviewed By :Jagrut Upadhyay 01/19/2022
 Supervised By :mohammad ahmed 01/21/2022



TIC: BG052057.D\data.ms

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

3.572min (-0.005) 5.00 ng/uL

response 3170

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	85.47
34.00	0.00	0.00
0.00	0.00	0.00

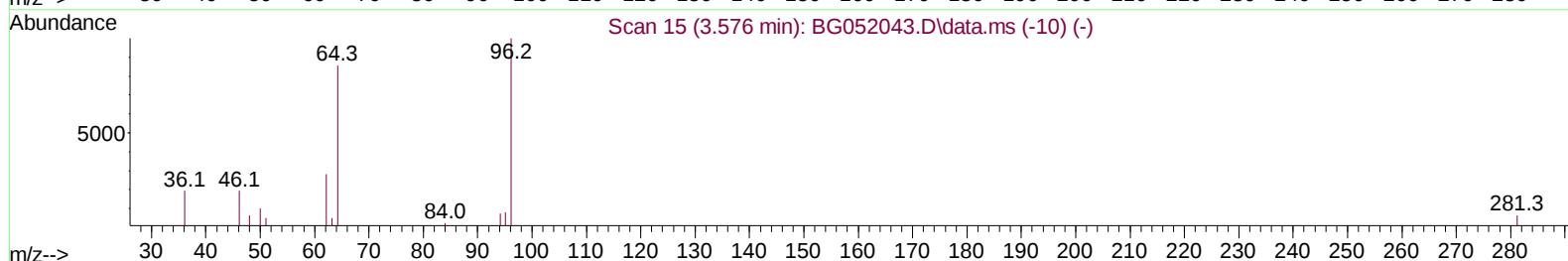
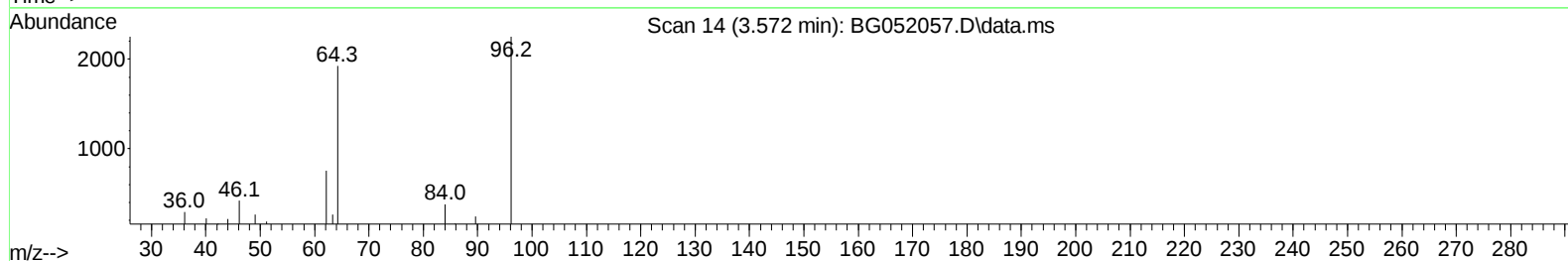
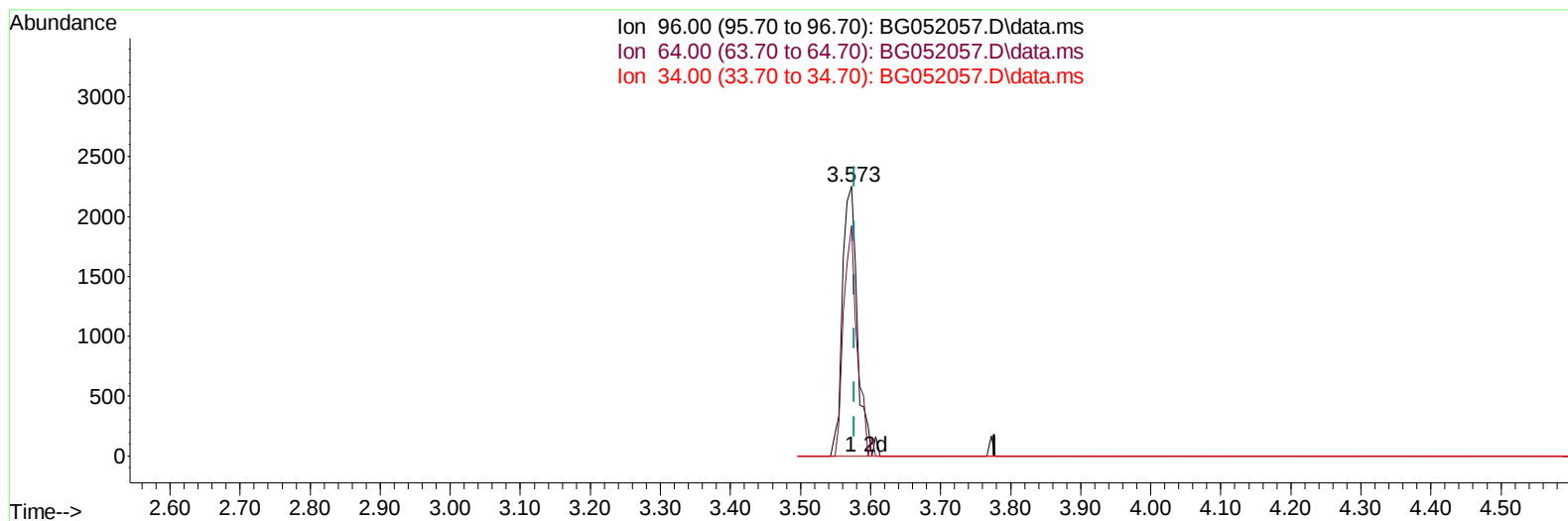
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TIC: BG052057.D\data.ms

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

3.572min (-0.005) 5.14 ng/uL m

response 3259

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	77.60	85.47
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	8.243	152	22668	20.000	ng/ul	-0.01
20) Naphthalene-d8	11.086	136	103588	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.887	164	77636	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.636	188	186649	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.953	240	171118	20.000	ng/ul	# 0.00
88) Perylene-d12	25.437	264	171512	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.572	96	3259m	5.137	ng/uL	0.00
4) Pyridine-d5	3.995	84	50053	25.382	ng/ul	-0.01
7) Phenol-d5	7.385	99	62971	28.383	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.544	67	37889	26.903	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.767	132	44916	30.578	ng/ul	-0.01
15) 4-Methylphenol-d8	8.942	113	51616	29.928	ng/ul	-0.01
21) Nitrobenzene-d5	9.412	128	22917	27.998	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.146	143	29183	32.237	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.692	165	52798	29.838	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.209	131	70533	28.623	ng/ul	0.00
46) Dimethylphthalate-d6	14.276	166	184282	28.717	ng/ul	0.00
49) Acenaphthylene-d8	14.581	160	225946	29.148	ng/ul	0.00
54) 4-Nitrophenol-d4	15.063	143	26430	24.634	ng/ul	0.00
60) Fluorene-d10	15.874	176	160023	28.601	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.985	200	25693	21.821	ng/ul	0.00
73) Anthracene-d10	17.736	188	276173	31.191	ng/ul	0.00
81) Pyrene-d10	20.015	212	319890	32.482	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.202	264	284955	31.559	ng/ul	0.00

Target Compounds Qvalue

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

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