

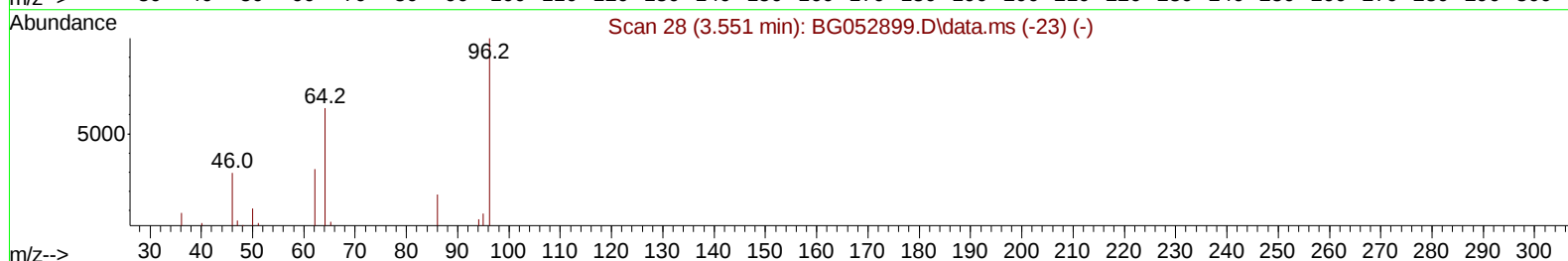
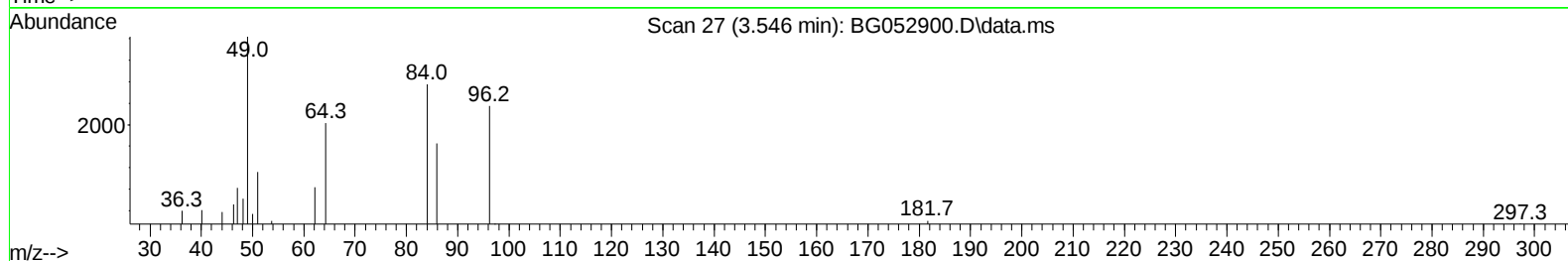
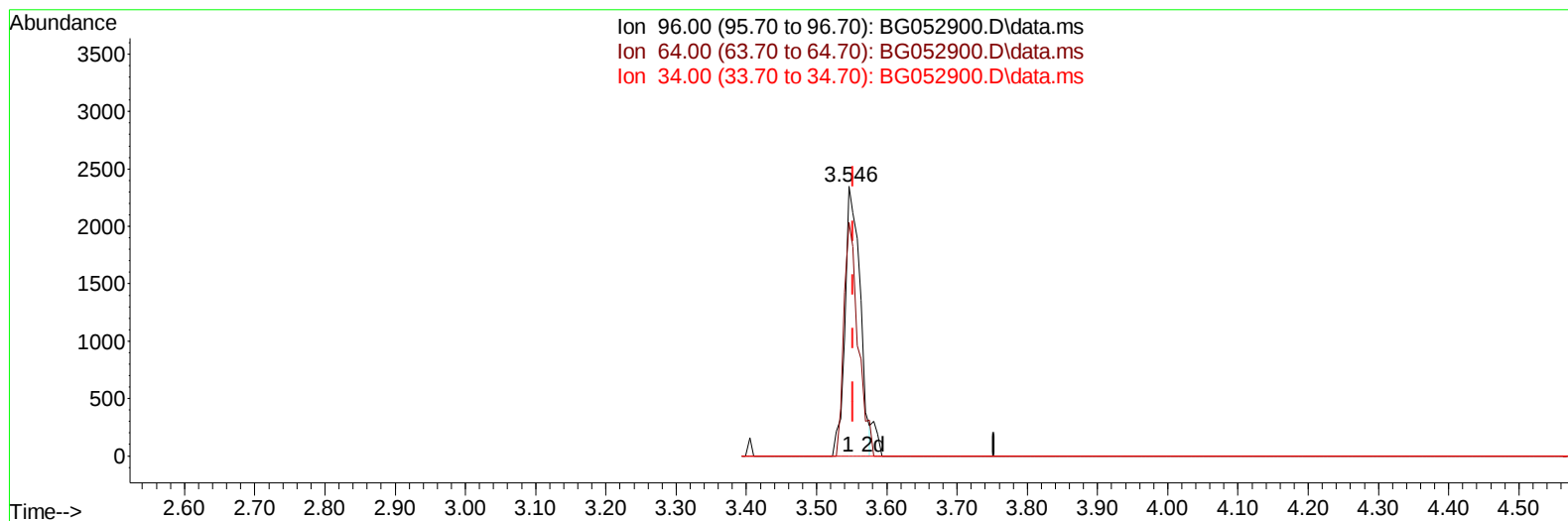
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG032822\
 Data File : BG052900.D
 Acq On : 28 Mar 2022 21:14
 Operator : CG/JU
 Sample : PB143650BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK650

Manual IntegrationsAPPROVED

Quant Time: Mar 29 02:35:53 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 03/29/2022
 Supervised By :mohammad ahmed 03/29/2022



TIC: BG052900.D\data.ms

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

3.546min (-0.007) 5.04 ng/uL

response 3499

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	86.66
34.00	0.00	0.00
0.00	0.00	0.00

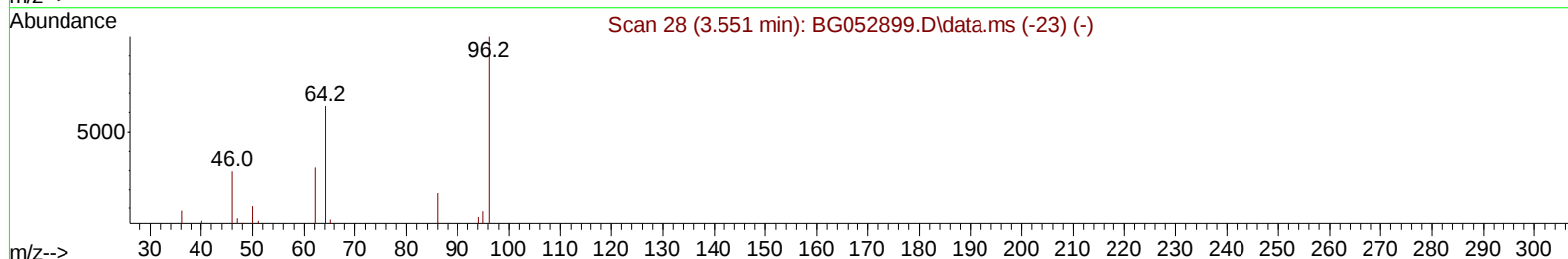
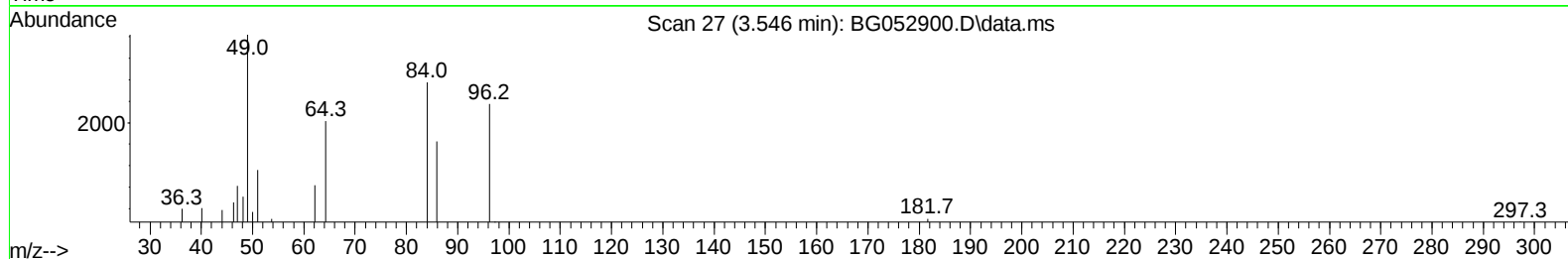
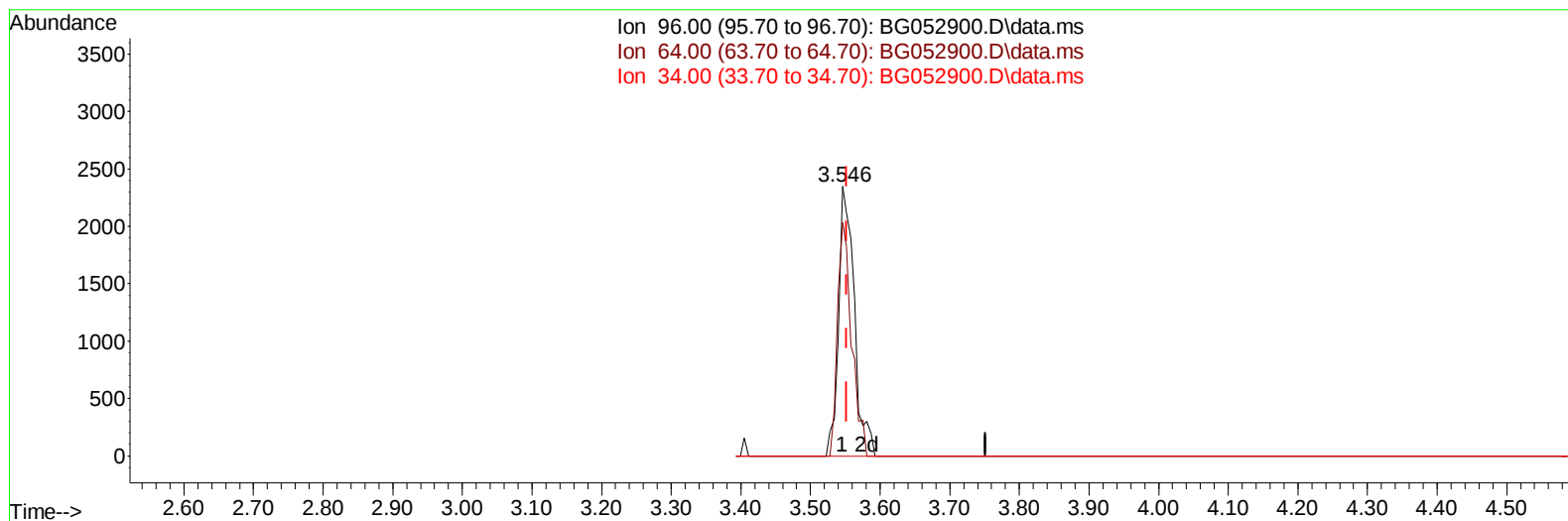
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TIC: BG052900.D\data.ms

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

3.546min (-0.007) 5.29 ng/uL m

response 3674

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	83.40	86.66
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	8.140	152	24538	20.000	ng/ul	0.00
20) Naphthalene-d8	10.948	136	113607	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.754	164	94336	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.498	188	227806	20.000	ng/ul	0.00
79) Chrysene-d12	21.786	240	256919	20.000	ng/ul	# 0.00
88) Perylene-d12	25.099	264	251407	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.546	96	3674m	5.289	ng/uL	0.00
4) Pyridine-d5	3.963	84	46482	26.802	ng/ul	-0.01
7) Phenol-d5	7.282	99	65520	29.812	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.452	67	40824	30.414	ng/ul	0.00
11) 2-Chlorophenol-d4	7.664	132	44073	29.534	ng/ul	-0.01
15) 4-Methylphenol-d8	8.827	113	49589	31.148	ng/ul	0.00
21) Nitrobenzene-d5	9.291	128	24889	30.435	ng/ul	0.00
24) 2-Nitrophenol-d4	10.019	143	29237	33.896	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.560	165	52204	27.820	ng/ul	0.00
31) 4-Chloroaniline-d4	11.071	131	74296	29.957	ng/ul	0.00
46) Dimethylphthalate-d6	14.149	166	227357	31.371	ng/ul	-0.01
49) Acenaphthylene-d8	14.449	160	244342	27.746	ng/ul	0.00
54) 4-Nitrophenol-d4	14.925	143	36390	29.865	ng/ul	0.00
60) Fluorene-d10	15.741	176	195861	30.779	ng/ul	-0.01
65) 4,6-Dinitro-2-methylph...	15.847	200	34150	29.457	ng/ul	-0.01
73) Anthracene-d10	17.597	188	348357	32.839	ng/ul	0.00
81) Pyrene-d10	19.877	212	444023	30.894	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.870	264	435852	33.552	ng/ul	0.00

Target Compounds Qvalue

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

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