

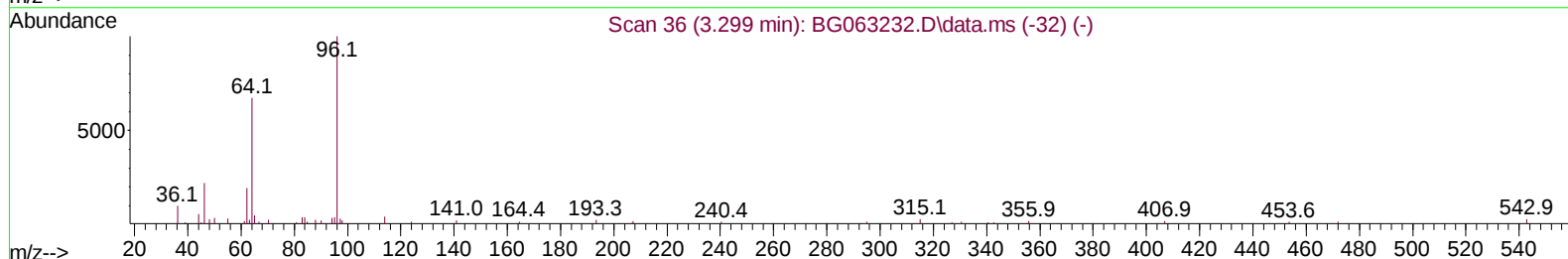
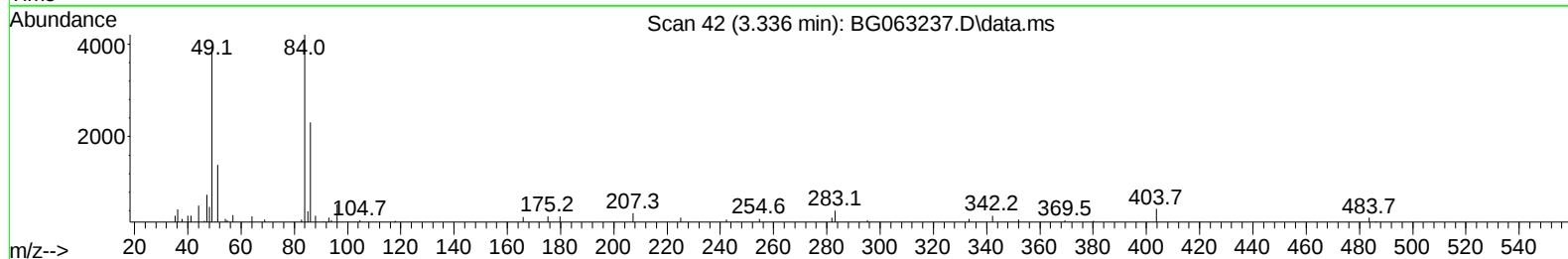
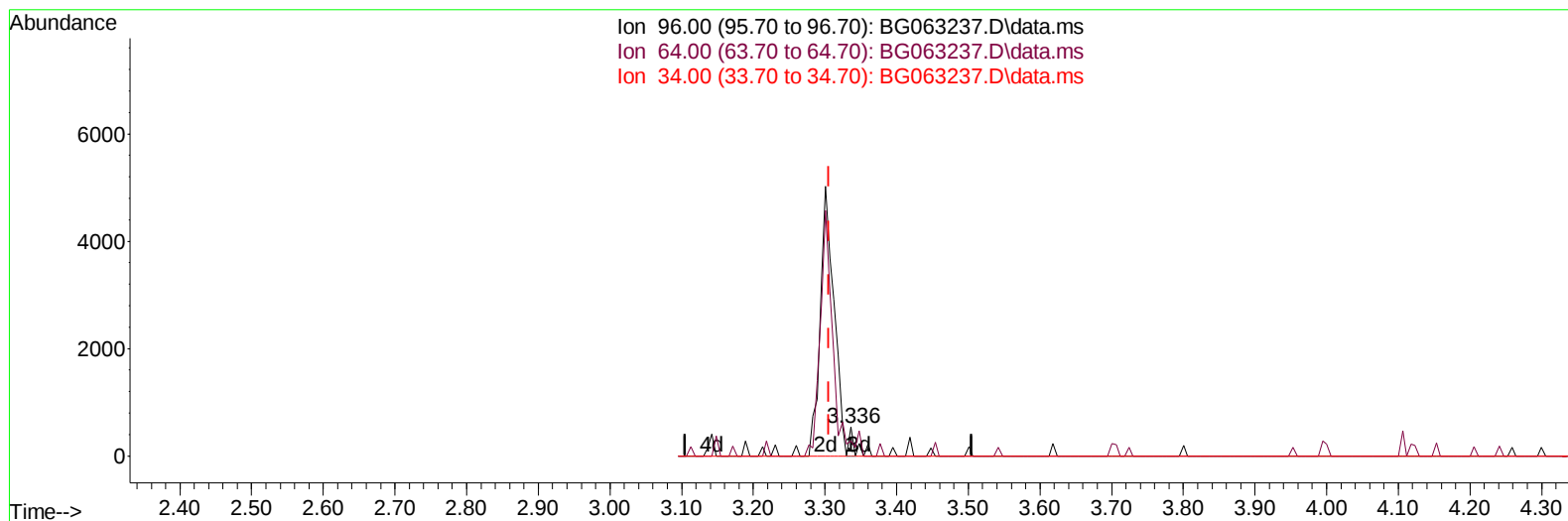
Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG110824\
Data File : BG063237.D
Acq On : 8 Nov 2024 20:31
Operator : RC/JU
Sample : P4636-10
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_G
ClientSampleId :
CC0Q2

Manual IntegrationsAPPROVED

Quant Time: Nov 08 22:12:35 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110624.MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Nov 06 15:45:52 2024
Response via : Initial Calibration

Reviewed By :Yogesh Patel 11/10/2024
Supervised By :mohammad ahmed 11/11/2024



TIC: BG063237.D\data.ms

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

3.336min (+ 0.031) 0.09 ng/uL

response 188

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	52.20	52.06
34.00	0.00	0.00
0.00	0.00	0.00

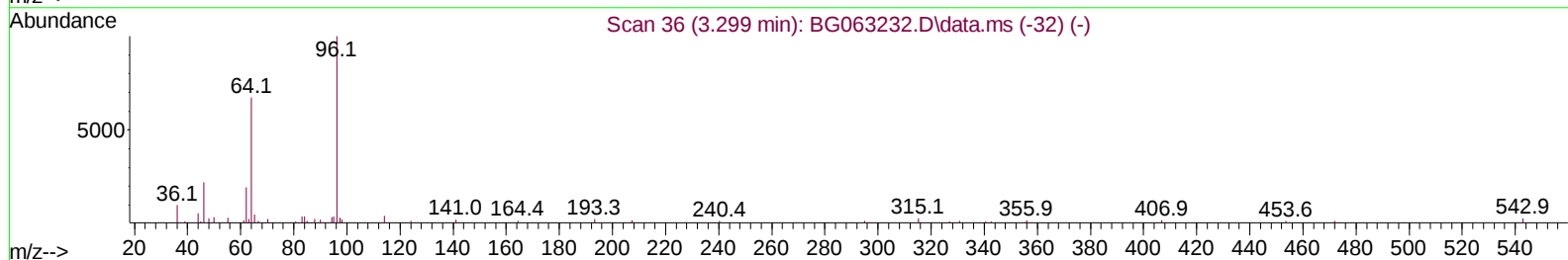
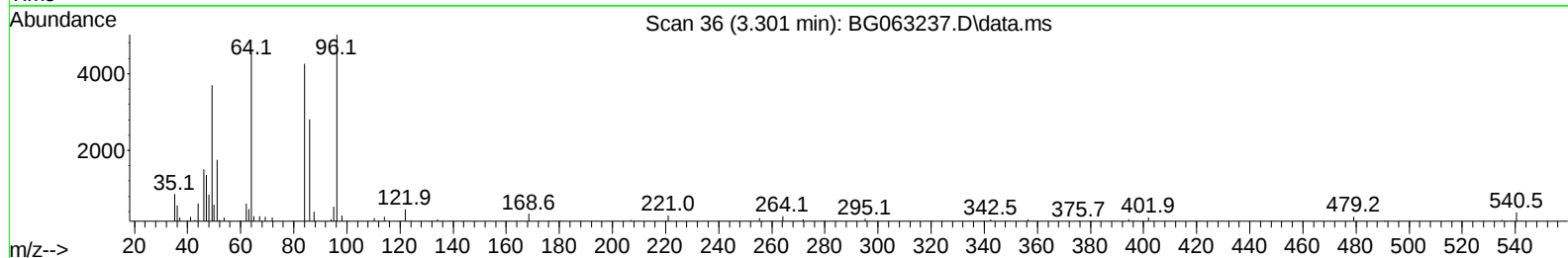
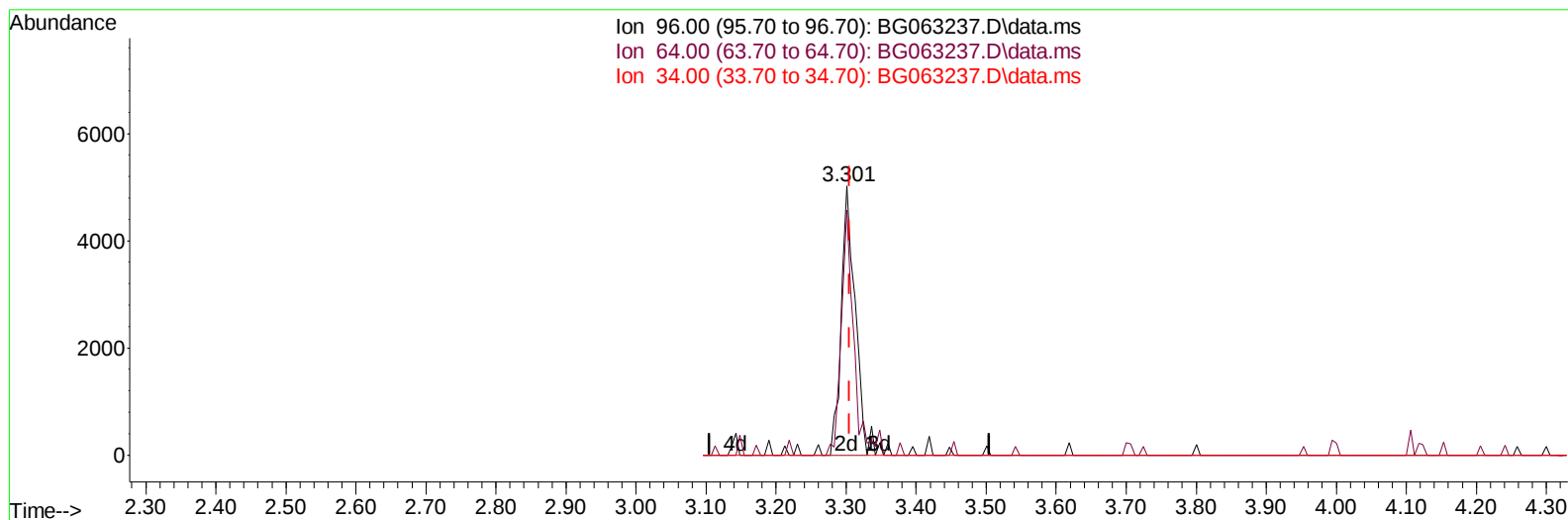
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(3) 1,4-Dioxane-d8 (S)

3.301min (-0.004) 3.24 ng/uL m

response 6774

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	52.20	91.20#
34.00	0.00	0.00
0.00	0.00	0.00

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 QLast Update : Wed Nov 06 15:45:52 2024
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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.837	152		94564	20.000	ng/ul	0.00
20) Naphthalene-d8	10.628	136		435954	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.470	164		389468	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.226	188		954318	20.000	ng/ul	0.00
79) Chrysene-d12	21.480	240		953456	20.000	ng/ul	# 0.00
88) Perylene-d12	24.511	264		1018147	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.301	96		6774m	3.239	ng/uL	0.00
4) Pyridine-d5	3.706	84		81482	11.788	ng/ul	0.00
7) Phenol-d5	7.008	99		151719	17.780	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.167	67		88303	17.643	ng/ul	0.00
11) 2-Chlorophenol-d4	7.367	132		104978	18.954	ng/ul	-0.01
15) 4-Methylphenol-d8	8.542	113		128384	17.751	ng/ul	0.00
21) Nitrobenzene-d5	8.988	128		59876	19.536	ng/ul	0.00
24) 2-Nitrophenol-d4	9.711	143		77616	19.594	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.252	165		150705	19.386	ng/ul	0.00
31) 4-Chloroaniline-d4	10.763	131		156342	15.441	ng/ul	0.00
46) Dimethylphthalate-d6	13.877	166		561003	18.389	ng/ul	0.00
49) Acenaphthylene-d8	14.165	160		614463	20.578	ng/ul	0.00
54) 4-Nitrophenol-d4	14.682	143		65520	15.686	ng/ul	0.00
60) Fluorene-d10	15.469	176		522685	20.301	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.592	200		95187	15.871	ng/ul	0.00
73) Anthracene-d10	17.320	188		830625	20.252	ng/ul	0.00
81) Pyrene-d10	19.623	212		1088050	22.761	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.306	264		1054207	21.446	ng/ul	0.00

Target Compounds Qvalue

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

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