

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG092324\
 Data File : BG063012.D
 Acq On : 24 Sep 2024 4:21
 Operator : RC/JU
 Sample : PB163509BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

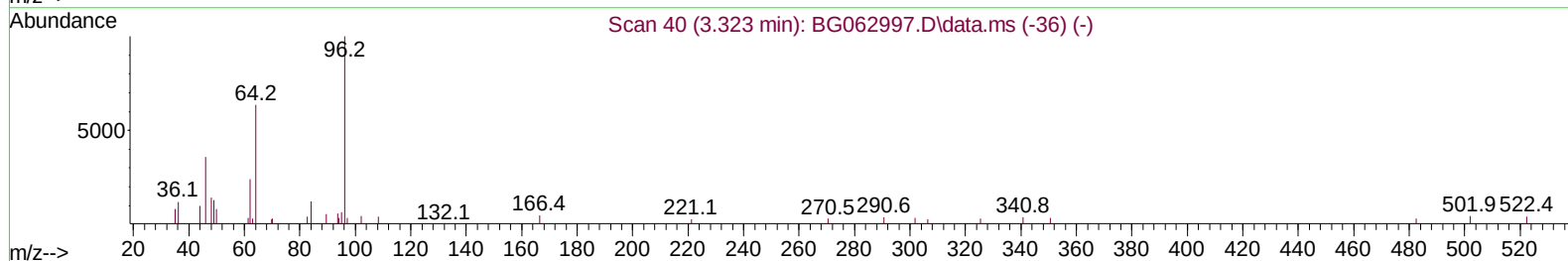
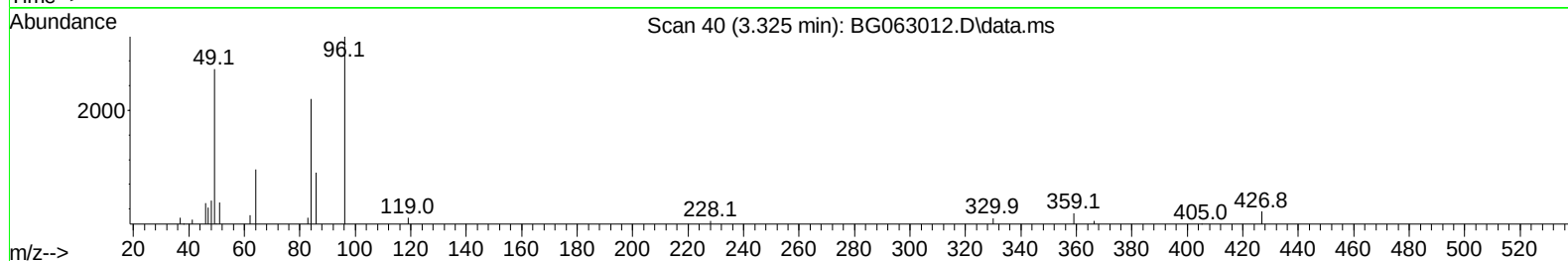
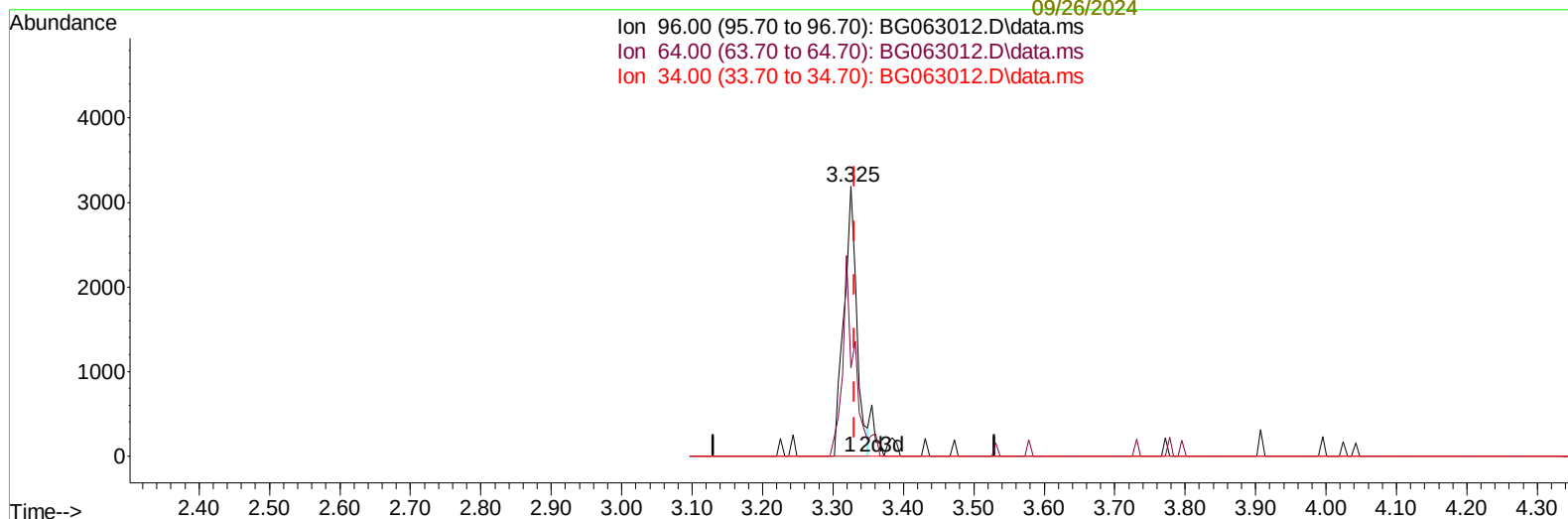
Instrument :
 BNA_G
 ClientSampleId :
 SBLK509

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 09/25/2024
 Supervised By :mohammad ahmed 09/26/2024

Supervised By :mohammad
 ahmed
 09/26/2024

Quant Time: Sep 24 04:56:53 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091724.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Sep 18 00:45:18 2024
 Response via : Initial Calibration



TIC: BG063012.D\data.ms

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

3.325min (-0.005) 4.95 ng/uL

response 3952

Ion	Exp%	Act%
96.00	100.00	100.00
64.00	75.20	32.67#
34.00	0.00	0.00
0.00	0.00	0.00

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Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)

Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.850	152		23392	20.000	ng/ul	0.00
20) Naphthalene-d8	10.640	136		101604	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.489	164		89570	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.233	188		237978	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.486	240		265936	20.000	ng/ul	# 0.00
88) Perylene-d12	24.530	264		327244	20.000	ng/ul	0.01
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.325	96		4277m	5.355	ng/uL	0.00
4) Pyridine-d5	3.725	84		39486	18.212	ng/ul	-0.01
7) Phenol-d5	7.027	99		75522	31.114	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.186	67		40359	29.712	ng/ul	0.00
11) 2-Chlorophenol-d4	7.385	132		48282	30.297	ng/ul	0.00
15) 4-Methylphenol-d8	8.560	113		58948	33.152	ng/ul	0.00
21) Nitrobenzene-d5	9.007	128		25930	32.034	ng/ul	0.00
24) 2-Nitrophenol-d4	9.730	143		35444	38.597	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.270	165		66020	34.651	ng/ul	0.00
31) 4-Chloroaniline-d4	10.776	131		78459	33.090	ng/ul	0.00
46) Dimethylphthalate-d6	13.895	166		253566	38.069	ng/ul	0.00
49) Acenaphthylene-d8	14.177	160		252403	33.928	ng/ul	0.00
54) 4-Nitrophenol-d4	14.694	143		34002	31.856	ng/ul	0.00
60) Fluorene-d10	15.482	176		229438	36.995	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.605	200		54388	37.807	ng/ul	0.00
73) Anthracene-d10	17.333	188		396091	35.217	ng/ul	0.00
81) Pyrene-d10	19.636	212		538496	33.433	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.324	264		581917	33.718	ng/ul	0.01

Target Compounds Qvalue

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

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