

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111324\
 Data File : BG063410.D
 Acq On : 15 Nov 2024 3:32
 Operator : RC/JU
 Sample : P4715-15
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0AL3

Manual IntegrationsAPPROVED

Reviewed By :Yogesh Patel 11/15/2024
 Supervised By :mohammad ahmed 11/18/2024

Quant Time: Nov 15 03:07:07 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

11/15/2024
 Supervised By :mohammad
 ahmed

Compound	R.T.	Q	Ion	Response	Conc	Units	Dev(Min)
-----11/18/2024							
Internal Standards							
1) 1,4-Dichlorobenzene-d4	7.832	152		115431	20.000	ng/ul	0.00
20) Naphthalene-d8	10.623	136		470936	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.471	164		389559	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.221	188		980772	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.475	240		993045	20.000	ng/ul	# 0.00
88) Perylene-d12	24.512	264		1068248	20.000	ng/ul	0.00
System Monitoring Compounds							
3) 1,4-Dioxane-d8	3.308	96		11272	4.416	ng/uL	0.00
4) Pyridine-d5	0.000	84		0d	0.000	ng/ul	
7) Phenol-d5	7.009	99		73511	7.058	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.162	67		136081	22.274	ng/ul	0.00
11) 2-Chlorophenol-d4	7.368	132		151009	22.336	ng/ul	0.00
15) 4-Methylphenol-d8	8.543	113		140643	15.931	ng/ul	0.00
21) Nitrobenzene-d5	8.989	128		81419	24.592	ng/ul	0.00
24) 2-Nitrophenol-d4	9.706	143		101608	23.745	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.247	165		200272	23.849	ng/ul	0.00
31) 4-Chloroaniline-d4	10.758	131		205652	18.802	ng/ul	0.00
46) Dimethylphthalate-d6	13.878	166		697011	22.842	ng/ul	0.00
49) Acenaphthylene-d8	14.160	160		708648	23.727	ng/ul	0.00
54) 4-Nitrophenol-d4	14.694	143		27393m	6.557	ng/ul	0.01
60) Fluorene-d10	15.464	176		627235	24.356	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.593	200		143609	23.298	ng/ul	0.00
73) Anthracene-d10	17.321	188		1106350	26.247	ng/ul	0.00
81) Pyrene-d10	19.618	212		1404451	28.208	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.301	264		1356469	26.300	ng/ul	0.00

Target Compounds Qvalue

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

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