

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122624\
 Data File : BG063872.D
 Acq On : 27 Dec 2024 17:25
 Operator : RC/JU
 Sample : P5323-01
 Misc : GCMS CONFIRMATION
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 YE5X0

Quant Time: Dec 27 23:31:48 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121124.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 24 00:44:39 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.794	152	109508	20.000	ng/ul	0.00
20) Naphthalene-d8	10.585	136	415274	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.439	164	263922	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.195	188	629607	20.000	ng/ul	0.00
79) Chrysene-d12	21.443	240	766175	20.000	ng/ul	0.00
88) Perylene-d12	24.463	264	780371	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.252	96	57	0.020	ng/uL	0.00
4) Pyridine-d5	3.658	84	211	0.024	ng/ul	0.00
7) Phenol-d5	0.000	99	0	0.000	ng/ul	
9) Bis-(2-Chloroethyl)eth...	7.142	67	79	0.011	ng/ul	0.01
11) 2-Chlorophenol-d4	7.159	132	59	0.009	ng/ul	-0.18
15) 4-Methylphenol-d8	0.000	113	0	0.000	ng/ul	
21) Nitrobenzene-d5	0.000	128	0	0.000	ng/ul	
24) 2-Nitrophenol-d4	0.000	143	0	0.000	ng/ul	
28) 2,4-Dichlorophenol-d3	0.000	165	0	0.000	ng/ul	
31) 4-Chloroaniline-d4	10.614	131	241	0.029	ng/ul	-0.11
46) Dimethylphthalate-d6	0.000	166	0	0.000	ng/ul	
49) Acenaphthylene-d8	0.000	160	0	0.000	ng/ul	
54) 4-Nitrophenol-d4	14.604	143	80	0.031	ng/ul	-0.06
60) Fluorene-d10	15.450	176	65	0.004	ng/ul	0.02
65) 4,6-Dinitro-2-methylph...	15.526	200	113	0.035	ng/ul	-0.04
73) Anthracene-d10	17.195	188	629607	23.259	ng/ul	-0.09
81) Pyrene-d10	19.657	212	83	0.002	ng/ul	0.07
92) Benzo(a)pyrene-d12	0.000	264	0	0.000	ng/ul	

Target Compounds Qvalue

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

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