

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122624\
 Data File : BG063870.D
 Acq On : 27 Dec 2024 16:10
 Operator : RC/JU
 Sample : PB165816BL
 Misc : GCMS CONFIRMATION
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK816

Quant Time: Dec 27 23:31:24 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	7.792	152	105964	20.000	ng/ul	0.00
20) Naphthalene-d8	10.582	136	401552	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.443	164	254745	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.192	188	614318	20.000	ng/ul	0.00
79) Chrysene-d12	21.446	240	664872	20.000	ng/ul	0.00
88) Perylene-d12	24.460	264	680539	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.000	96	0	0.000	ng/uL	
4) Pyridine-d5	3.649	84	295	0.035	ng/ul	-0.01
7) Phenol-d5	0.000	99	0	0.000	ng/ul	
9) Bis-(2-Chloroethyl)eth...	6.998	67	191	0.028	ng/ul	-0.13
11) 2-Chlorophenol-d4	0.000	132	0	0.000	ng/ul	
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	10.594	131	173	0.022	ng/ul	-0.13
46) Dimethylphthalate-d6	13.696	166	67	0.004	ng/ul	-0.15
49) Acenaphthylene-d8	0.000	160	0	0.000	ng/ul	
54) 4-Nitrophenol-d4	0.000	143	0	0.000	ng/ul	
60) Fluorene-d10	0.000	176	0	0.000	ng/ul	
65) 4,6-Dinitro-2-methylph...	0.000	200	0	0.000	ng/ul	
73) Anthracene-d10	17.192	188	614318	23.259	ng/ul	-0.10
81) Pyrene-d10	0.000	212	0	0.000	ng/ul	
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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