

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122624\
 Data File : BG063869.D
 Acq On : 27 Dec 2024 15:32
 Operator : RC/JU
 Sample : PB165831BL
 Misc : GCMS CONFIRMATION
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK831

Quant Time: Dec 27 23:31:16 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	98239	20.000	ng/ul	0.00
20) Naphthalene-d8	10.585	136	356115	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.439	164	233168	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.194	188	549845	20.000	ng/ul	0.00
79) Chrysene-d12	21.442	240	632509	20.000	ng/ul	0.00
88) Perylene-d12	24.462	264	643630	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.193	96	137	0.055	ng/uL	-0.06
4) Pyridine-d5	3.663	84	780	0.099	ng/ul	0.00
7) Phenol-d5	0.000	99	0	0.000	ng/ul	
9) Bis-(2-Chloroethyl)eth...	7.112	67	53	0.008	ng/ul	-0.02
11) 2-Chlorophenol-d4	7.418	132	208	0.036	ng/ul	0.08
15) 4-Methylphenol-d8	8.616	113	123	0.018	ng/ul	0.10
21) Nitrobenzene-d5	0.000	128	0	0.000	ng/ul	
24) 2-Nitrophenol-d4	9.715	143	64	0.022	ng/ul	0.04
28) 2,4-Dichlorophenol-d3	10.126	165	146	0.026	ng/ul	-0.09
31) 4-Chloroaniline-d4	10.802	131	82	0.012	ng/ul	0.08
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	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	17.194	188	549845	23.259	ng/ul	-0.09
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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