

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG091023\
 Data File : BG058977.D
 Acq On : 11 Sep 2023 10:40
 Operator : MA/JU
 Sample : PB155328BL
 Misc :
 ALS Vial : 76 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK328

Quant Time: Sep 12 01:06:29 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG091023.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Sep 11 00:07:05 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.329	152	56828	20.000	ng/u1	0.00
20) Naphthalene-d8	11.172	136	267624	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.950	164	199674	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.700	188	523987	20.000	ng/u1	# 0.00
79) Chrysene-d12	22.019	240	521056	20.000	ng/u1	0.00
88) Perylene-d12	25.591	264	611066	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.640	96	10287	7.694	ng/uL	0.00
4) Pyridine-d5	4.069	84	134532	31.782	ng/u1	0.00
7) Phenol-d5	7.436	99	172859	30.164	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.641	67	105833	31.495	ng/u1	0.00
11) 2-Chlorophenol-d4	7.841	132	112325	30.562	ng/u1	0.00
15) 4-Methylphenol-d8	9.004	113	136346	30.796	ng/u1	0.00
21) Nitrobenzene-d5	9.516	128	59513	30.657	ng/u1	0.00
24) 2-Nitrophenol-d4	10.238	143	70928	33.299	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.761	165	134841	29.786	ng/u1	0.00
31) 4-Chloroaniline-d4	11.308	131	181308	29.782	ng/u1	0.00
46) Dimethylphthalate-d6	14.345	166	499845	32.272	ng/u1	0.00
49) Acenaphthylene-d8	14.651	160	524407	30.405	ng/u1	0.00
54) 4-Nitrophenol-d4	15.115	143	85180	33.109	ng/u1	0.00
60) Fluorene-d10	15.938	176	437421	31.823	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	16.037	200	75799	29.470	ng/u1	0.00
73) Anthracene-d10	17.800	188	759897	32.668	ng/u1	0.00
81) Pyrene-d10	20.068	212	893143	32.358	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.338	264	941845	31.442	ng/u1	0.00

Target Compounds Qvalue

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

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