

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010622\
 Data File : BG051893.D
 Acq On : 6 Jan 2022 15:01
 Operator : CG/JU
 Sample : PB141855BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK855

Quant Time: Jan 06 23:56:22 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 06 14:43:02 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.252	152	25875	20.000	ng/u1	0.00
20) Naphthalene-d8	11.090	136	112237	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.896	164	83435	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.640	188	211513	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.957	240	231997	20.000	ng/u1	# 0.00
88) Perylene-d12	25.453	264	222257	20.000	ng/u1	0.01

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.576	96	4914	6.786	ng/uL	0.00
4) Pyridine-d5	3.999	84	67076	29.798	ng/u1	0.00
7) Phenol-d5	7.389	99	79255	31.296	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.559	67	53150	33.061	ng/u1	0.00
11) 2-Chlorophenol-d4	7.776	132	55633	33.180	ng/u1	0.00
15) 4-Methylphenol-d8	8.951	113	60778	30.873	ng/u1	0.00
21) Nitrobenzene-d5	9.421	128	30030	33.861	ng/u1	0.00
24) 2-Nitrophenol-d4	10.156	143	33027	33.672	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.702	165	59957	31.273	ng/u1	0.00
31) 4-Chloroaniline-d4	11.219	131	80681	30.218	ng/u1	0.00
46) Dimethylphthalate-d6	14.285	166	232243	33.675	ng/u1	0.00
49) Acenaphthylene-d8	14.591	160	266443	31.983	ng/u1	0.00
54) 4-Nitrophenol-d4	15.067	143	35694	30.956	ng/u1	0.00
60) Fluorene-d10	15.883	176	193361	32.157	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.989	200	38736	29.030	ng/u1	0.00
73) Anthracene-d10	17.739	188	340360	33.922	ng/u1	0.00
81) Pyrene-d10	20.019	212	437753	32.786	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.212	264	408732	34.933	ng/u1	0.02

Target Compounds	Qvalue

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

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