

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122021\
 Data File : BG051676.D
 Acq On : 20 Dec 2021 11:16
 Operator : CG/JU
 Sample : PB141560BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK560

Quant Time: Dec 20 23:19:28 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Tue Dec 14 14:19:57 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.274	152	37274	20.000	ng/u1	-0.01
20) Naphthalene-d8	11.112	136	155022	20.000	ng/u1	#-0.01
38) Acenaphthene-d10	14.907	164	111413	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.650	188	256097	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.974	240	252021	20.000	ng/u1	#-0.01
88) Perylene-d12	25.463	264	245612	20.000	ng/u1	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.598	96	5417	6.036	ng/uL	0.00
4) Pyridine-d5	4.021	84	74854	27.603	ng/u1	-0.02
7) Phenol-d5	7.405	99	105889	31.817	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.575	67	62378	31.528	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.793	132	73953	31.950	ng/u1	-0.02
15) 4-Methylphenol-d8	8.968	113	82461	32.054	ng/u1	-0.01
21) Nitrobenzene-d5	9.443	128	37505	32.999	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.178	143	41778	33.625	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.718	165	82522	30.628	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.229	131	89423	25.208	ng/u1	-0.01
46) Dimethylphthalate-d6	14.296	166	272417	30.549	ng/u1	-0.02
49) Acenaphthylene-d8	14.601	160	336349	30.370	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.065	143	37249	25.686	ng/u1	-0.01
60) Fluorene-d10	15.894	176	233590	29.237	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.993	200	32011	23.991	ng/u1	-0.02
73) Anthracene-d10	17.756	188	390717	31.983	ng/u1	0.00
81) Pyrene-d10	20.029	212	462441	30.488	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.228	264	421972	32.652	ng/u1	-0.02

Target Compounds Qvalue

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

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