

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
 Data File : BG051785.D
 Acq On : 23 Dec 2021 15:02
 Operator : CG/JU
 Sample : PB141646BL
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK646

Quant Time: Dec 24 00:11:36 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Dec 23 13:37:07 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.271	152	35507	20.000	ng/u1	-0.02
20) Naphthalene-d8	11.108	136	144299	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.909	164	102234	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.652	188	232246	20.000	ng/u1	#-0.01
79) Chrysene-d12	21.970	240	228358	20.000	ng/u1	#-0.02
88) Perylene-d12	25.465	264	217752	20.000	ng/u1	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.600	96	5217	6.102	ng/uL	0.00
4) Pyridine-d5	4.023	84	83779	32.432	ng/u1	-0.02
7) Phenol-d5	7.401	99	101087	31.885	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.572	67	61109	32.424	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.795	132	70932	32.170	ng/u1	-0.02
15) 4-Methylphenol-d8	8.964	113	82133	33.516	ng/u1	-0.02
21) Nitrobenzene-d5	9.440	128	38205	36.113	ng/u1	-0.02
24) 2-Nitrophenol-d4	10.174	143	41611	35.980	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.714	165	82242	32.792	ng/u1	-0.02
31) 4-Chloroaniline-d4	11.231	131	104746	31.722	ng/u1	-0.01
46) Dimethylphthalate-d6	14.292	166	262173	32.040	ng/u1	-0.02
49) Acenaphthylene-d8	14.603	160	323398	31.822	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.073	143	35640	26.783	ng/u1	0.00
60) Fluorene-d10	15.896	176	226927	30.954	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.995	200	30407	25.129	ng/u1	-0.02
73) Anthracene-d10	17.752	188	370733	33.463	ng/u1	-0.01
81) Pyrene-d10	20.025	212	445441	32.410	ng/u1	-0.02
92) Benzo(a)pyrene-d12	25.224	264	402618	35.141	ng/u1	-0.03

Target Compounds	Qvalue

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

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