

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
 Data File : BG051635.D
 Acq On : 18 Dec 2021 18:43
 Operator : CG/JU
 Sample : PB141535BL
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK535

Quant Time: Dec 18 20:53:15 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	8.273	152	38305	20.000	ng/u1	-0.01
20) Naphthalene-d8	11.110	136	161449	20.000	ng/u1	#-0.01
38) Acenaphthene-d10	14.911	164	117155	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.654	188	273123	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.978	240	271878	20.000	ng/u1	# 0.00
88) Perylene-d12	25.473	264	262052	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.603	96	5500	5.963	ng/uL	0.00
4) Pyridine-d5	4.026	84	81937	29.402	ng/u1	-0.02
7) Phenol-d5	7.403	99	103463	30.251	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.580	67	62337	30.660	ng/u1	-0.01
11) 2-Chlorophenol-d4	7.797	132	74694	31.401	ng/u1	-0.01
15) 4-Methylphenol-d8	8.966	113	81160	30.699	ng/u1	-0.01
21) Nitrobenzene-d5	9.442	128	38403	32.444	ng/u1	-0.01
24) 2-Nitrophenol-d4	10.176	143	42720	33.015	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.717	165	80839	28.809	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.228	131	109572	29.658	ng/u1	-0.01
46) Dimethylphthalate-d6	14.300	166	279742	29.833	ng/u1	-0.01
49) Acenaphthylene-d8	14.605	160	340161	29.209	ng/u1	0.00
54) 4-Nitrophenol-d4	15.069	143	41023	26.902	ng/u1	0.00
60) Fluorene-d10	15.898	176	240359	28.610	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.998	200	36944	25.962	ng/u1	-0.01
73) Anthracene-d10	17.754	188	398181	30.562	ng/u1	0.00
81) Pyrene-d10	20.027	212	472715	28.889	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.232	264	431847	31.320	ng/u1	-0.02

Target Compounds Qvalue

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

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