

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122221\
 Data File : BG051767.D
 Acq On : 23 Dec 2021 2:43
 Operator : CG/JU
 Sample : PB141620BL
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK620

Quant Time: Dec 23 05:52:15 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	8.272	152	41534	20.000	ng/u1	-0.02
20) Naphthalene-d8	11.109	136	175105	20.000	ng/u1	-0.02
38) Acenaphthene-d10	14.910	164	127633	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.653	188	291405	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.977	240	286806	20.000	ng/u1	# 0.00
88) Perylene-d12	25.472	264	273158	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.596	96	5399	5.398	ng/uL	0.00
4) Pyridine-d5	4.024	84	76907	25.452	ng/u1	-0.02
7) Phenol-d5	7.402	99	113427	30.586	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.573	67	66385	30.112	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.796	132	78641	30.491	ng/u1	-0.02
15) 4-Methylphenol-d8	8.965	113	91013	31.750	ng/u1	-0.02
21) Nitrobenzene-d5	9.441	128	40367	31.443	ng/u1	-0.02
24) 2-Nitrophenol-d4	10.169	143	47163	33.606	ng/u1	-0.02
28) 2,4-Dichlorophenol-d3	10.715	165	88094	28.946	ng/u1	-0.02
31) 4-Chloroaniline-d4	11.232	131	90001	22.461	ng/u1	0.00
46) Dimethylphthalate-d6	14.293	166	291801	28.564	ng/u1	-0.02
49) Acenaphthylene-d8	14.604	160	365818	28.833	ng/u1	0.00
54) 4-Nitrophenol-d4	15.074	143	39392	23.712	ng/u1	0.00
60) Fluorene-d10	15.897	176	255990	27.969	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.996	200	37452	24.668	ng/u1	-0.02
73) Anthracene-d10	17.753	188	413490	29.746	ng/u1	0.00
81) Pyrene-d10	20.032	212	496554	28.767	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.225	264	448296	31.191	ng/u1	-0.03

Target Compounds Qvalue

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

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