

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG010322\
 Data File : BG051858.D
 Acq On : 3 Jan 2022 15:29
 Operator : CG/JU
 Sample : PB141780BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK780

Quant Time: Jan 04 01:22:29 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG121421.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Fri Dec 31 00:13:36 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.271	152	25380	20.000	ng/u1	0.00
20) Naphthalene-d8	11.109	136	98268	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.904	164	72037	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.647	188	193416	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.965	240	190424	20.000	ng/u1	# 0.00
88) Perylene-d12	25.460	264	176982	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.589	96	3789	6.200	ng/uL	0.00
4) Pyridine-d5	4.024	84	59519	32.234	ng/u1	0.00
7) Phenol-d5	7.408	99	67615	29.838	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.578	67	45317	33.639	ng/u1	0.00
11) 2-Chlorophenol-d4	7.796	132	47231	29.968	ng/u1	0.00
15) 4-Methylphenol-d8	8.970	113	52791	30.138	ng/u1	0.00
21) Nitrobenzene-d5	9.440	128	23845	33.097	ng/u1	0.00
24) 2-Nitrophenol-d4	10.169	143	25921	32.912	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.721	165	49114	28.756	ng/u1	0.00
31) 4-Chloroaniline-d4	11.232	131	68331	30.387	ng/u1	0.00
46) Dimethylphthalate-d6	14.293	166	193563	33.571	ng/u1	0.00
49) Acenaphthylene-d8	14.598	160	226228	31.592	ng/u1	0.00
54) 4-Nitrophenol-d4	15.074	143	31181	33.255	ng/u1	0.00
60) Fluorene-d10	15.890	176	167389	32.403	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.996	200	29413	29.188	ng/u1	0.00
73) Anthracene-d10	17.747	188	308013	33.383	ng/u1	0.00
81) Pyrene-d10	20.026	212	387689	33.828	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.219	264	306968	32.964	ng/u1	0.00

Target Compounds Qvalue

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

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