

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG012622\
 Data File : BG052206.D
 Acq On : 28 Jan 2022 4:39
 Operator : CG/JU
 Sample : SP5810
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SP5810

Quant Time: Jan 28 05:18:52 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Jan 26 23:58:52 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.242	152	25434	20.000	ng/u1	0.00
20) Naphthalene-d8	11.073	136	103095	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.885	164	73262	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.629	188	180456	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.952	240	185479	20.000	ng/u1	# 0.00
88) Perylene-d12	25.430	264	185108	20.000	ng/u1	0.00

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.571	96	11795	16.571	ng/uL	0.00
4) Pyridine-d5	3.988	84	171145	77.349	ng/u1	0.00
7) Phenol-d5	7.384	99	194092	77.970	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.542	67	117959	74.647	ng/u1	0.00
11) 2-Chlorophenol-d4	7.766	132	138527	84.051	ng/u1	0.00
15) 4-Methylphenol-d8	8.941	113	154784	79.988	ng/u1	0.00
21) Nitrobenzene-d5	9.411	128	73365	90.059	ng/u1	0.00
24) 2-Nitrophenol-d4	10.145	143	84096	93.341	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.691	165	156882	89.084	ng/u1	0.00
31) 4-Chloroaniline-d4	11.202	131	199794	81.466	ng/u1	0.00
46) Dimethylphthalate-d6	14.275	166	509187	84.084	ng/u1	0.00
49) Acenaphthylene-d8	14.580	160	640158	87.514	ng/u1	0.00
54) 4-Nitrophenol-d4	15.068	143	78907	77.936	ng/u1	0.00
60) Fluorene-d10	15.872	176	445199	84.321	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.984	200	81634	71.709	ng/u1	0.00
73) Anthracene-d10	17.729	188	738373	86.254	ng/u1	0.00
81) Pyrene-d10	20.014	212	893768	83.728	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.201	264	856969	87.940	ng/u1	0.02

Target Compounds	Qvalue

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

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