

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122021\
 Data File : BG051711.D
 Acq On : 21 Dec 2021 11:29
 Operator : CG/JU
 Sample : M5126-02
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0J11

Quant Time: Dec 21 12:06:00 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	8.278	152	38556	20.000	ng/ul	0.00
20) Naphthalene-d8	11.109	136	155590	20.000	ng/ul	#-0.01
38) Acenaphthene-d10	14.904	164	109054	20.000	ng/ul	-0.01
64) Phenanthrene-d10	17.653	188	257209	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.977	240	247076	20.000	ng/ul	# 0.00
88) Perylene-d12	25.472	264	247505	20.000	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.596	96	4070	4.384	ng/uL	0.00
4) Pyridine-d5	4.030	84	23965	8.544	ng/ul	-0.01
7) Phenol-d5	7.402	99	24011	6.975	ng/ul	-0.01
9) Bis-(2-Chloroethyl)eth...	7.579	67	60637	29.629	ng/ul	-0.01
11) 2-Chlorophenol-d4	7.796	132	53185	22.213	ng/ul	-0.01
15) 4-Methylphenol-d8	8.965	113	40877	15.361	ng/ul	-0.01
21) Nitrobenzene-d5	9.441	128	37196	32.607	ng/ul	-0.01
24) 2-Nitrophenol-d4	10.169	143	37318	29.926	ng/ul	-0.01
28) 2,4-Dichlorophenol-d3	10.716	165	75297	27.844	ng/ul	-0.01
31) 4-Chloroaniline-d4	11.227	131	86131	24.191	ng/ul	-0.01
46) Dimethylphthalate-d6	14.299	166	279650	32.038	ng/ul	-0.01
49) Acenaphthylene-d8	14.604	160	345394	31.861	ng/ul	0.00
54) 4-Nitrophenol-d4	15.068	143	8653	6.096	ng/ul	0.00
60) Fluorene-d10	15.897	176	250194	31.993	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.997	200	35895	26.786	ng/ul	-0.01
73) Anthracene-d10	17.753	188	423664	34.530	ng/ul	0.00
81) Pyrene-d10	20.032	212	489591	32.924	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.225	264	457914	35.163	ng/ul	-0.03

Target Compounds Qvalue

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

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