

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG111121\
 Data File : BG051018.D
 Acq On : 13 Nov 2021 4:58
 Operator : CG/JU
 Sample : M4618-03
 Misc :
 ALS Vial : 58 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BG206

Quant Time: Nov 15 01:27:01 2021
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG110321.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Mon Nov 15 00:27:19 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.237	152	46782	20.000	ng/u1	0.00
20) Naphthalene-d8	11.063	136	219484	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.859	164	142144	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.603	188	295069	20.000	ng/u1	-0.02
79) Chrysene-d12	21.904	240	244150	20.000	ng/u1	-0.02
88) Perylene-d12	25.323	264	235767	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.596	96	6115	4.219	ng/uL	0.00
4) Pyridine-d5	4.036	84	29255	6.747	ng/u1	-0.03
7) Phenol-d5	7.380	99	33827	6.778	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.550	67	99650	30.910	ng/u1	0.00
11) 2-Chlorophenol-d4	7.761	132	84214	24.348	ng/u1	0.00
15) 4-Methylphenol-d8	8.937	113	59964	15.262	ng/u1	0.00
21) Nitrobenzene-d5	9.412	128	62408	33.459	ng/u1	0.00
24) 2-Nitrophenol-d4	10.135	143	69131	33.333	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.676	165	101590	29.079	ng/u1	0.00
31) 4-Chloroaniline-d4	11.193	131	133236	25.184	ng/u1	-0.02
46) Dimethylphthalate-d6	14.254	166	410584	37.755	ng/u1	-0.02
49) Acenaphthylene-d8	14.559	160	499036	36.832	ng/u1	0.00
54) 4-Nitrophenol-d4	15.053	143	14567	7.388	ng/u1	0.00
60) Fluorene-d10	15.846	176	350538	36.387	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.964	200	60543	33.843	ng/u1	-0.02
73) Anthracene-d10	17.703	188	539285	38.657	ng/u1	0.00
81) Pyrene-d10	19.982	212	588566	37.324	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.082	264	495795	38.040	ng/u1	-0.02

Target Compounds	Qvalue

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

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