

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122021\
 Data File : BG051699.D
 Acq On : 21 Dec 2021 2:49
 Operator : CG/JU
 Sample : M5126-11
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C0J18

Quant Time: Dec 21 03:48:19 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	8.269	152	34726	20.000	ng/u1	-0.02
20) Naphthalene-d8	11.106	136	150744	20.000	ng/u1	#-0.02
38) Acenaphthene-d10	14.907	164	108395	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.656	188	246321	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.974	240	237970	20.000	ng/u1	#-0.01
88) Perylene-d12	25.469	264	230654	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.599	96	3685	4.407	ng/uL	0.00
4) Pyridine-d5	4.028	84	9807	3.882	ng/u1	-0.02
7) Phenol-d5	7.400	99	20346	6.562	ng/u1	-0.02
9) Bis-(2-Chloroethyl)eth...	7.576	67	57451	31.169	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.799	132	54087	25.082	ng/u1	-0.01
15) 4-Methylphenol-d8	8.962	113	37371	15.593	ng/u1	-0.02
21) Nitrobenzene-d5	9.438	128	36269	32.817	ng/u1	-0.02
24) 2-Nitrophenol-d4	10.172	143	37632	31.148	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.719	165	72755	27.769	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.230	131	71536	20.738	ng/u1	-0.01
46) Dimethylphthalate-d6	14.296	166	289215	33.335	ng/u1	-0.02
49) Acenaphthylene-d8	14.602	160	333111	30.915	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.072	143	7841	5.558	ng/u1	0.00
60) Fluorene-d10	15.894	176	251376	32.340	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.994	200	30557	23.810	ng/u1	-0.02
73) Anthracene-d10	17.756	188	433671	36.907	ng/u1	0.00
81) Pyrene-d10	20.030	212	515675	36.005	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.228	264	468948	38.641	ng/u1	-0.02

Target Compounds Qvalue

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

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