

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
 Data File : BG051624.D
 Acq On : 18 Dec 2021 10:34
 Operator : CG/JU
 Sample : M5037-21
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 C00Z0

Quant Time: Dec 18 20:02:40 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	8.276	152	30056	20.000	ng/u1	-0.01
20) Naphthalene-d8	11.113	136	124138	20.000	ng/u1	#-0.01
38) Acenaphthene-d10	14.908	164	87347	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.657	188	211900	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.975	240	202352	20.000	ng/u1	#-0.01
88) Perylene-d12	25.470	264	207959	20.000	ng/u1	-0.02

System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.599	96	3690	5.099	ng/uL	0.00
4) Pyridine-d5	4.028	84	22411	10.249	ng/u1	-0.02
7) Phenol-d5	7.406	99	9459	3.525	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.576	67	49241	30.866	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.800	132	25837	13.843	ng/u1	-0.01
15) 4-Methylphenol-d8	8.969	113	21329	10.282	ng/u1	-0.01
21) Nitrobenzene-d5	9.439	128	31151	34.227	ng/u1	-0.02
24) 2-Nitrophenol-d4	10.173	143	17729	17.819	ng/u1	-0.01
28) 2,4-Dichlorophenol-d3	10.719	165	34414	15.950	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.230	131	28957	10.194	ng/u1	-0.01
46) Dimethylphthalate-d6	14.297	166	248923	35.605	ng/u1	-0.02
49) Acenaphthylene-d8	14.602	160	277039	31.907	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.060	143	4128	3.631	ng/u1	-0.02
60) Fluorene-d10	15.900	176	223136	35.624	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	16.000	200	23967	21.709	ng/u1	-0.01
73) Anthracene-d10	17.757	188	405156	40.082	ng/u1	0.00
81) Pyrene-d10	20.030	212	473834	38.907	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.229	264	439137	40.133	ng/u1	-0.02

Target Compounds Qvalue

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

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