

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG121721\
 Data File : BG051647.D
 Acq On : 19 Dec 2021 2:57
 Operator : CG/JU
 Sample : M5154-13
 Misc :
 ALS Vial : 63 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGL93

Quant Time: Dec 19 03:30:33 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	38062	20.000	ng/u1	-0.01
20) Naphthalene-d8	11.113	136	167681	20.000	ng/u1	-0.01
38) Acenaphthene-d10	14.908	164	117844	20.000	ng/u1	-0.01
64) Phenanthrene-d10	17.657	188	279168	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.975	240	266822	20.000	ng/u1	#-0.01
88) Perylene-d12	25.470	264	268951	20.000	ng/u1	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.606	96	5521	6.024	ng/uL	0.00
4) Pyridine-d5	4.034	84	27994	10.109	ng/u1	-0.01
7) Phenol-d5	7.406	99	104763	30.827	ng/u1	-0.01
9) Bis-(2-Chloroethyl)eth...	7.577	67	60444	29.918	ng/u1	-0.02
11) 2-Chlorophenol-d4	7.800	132	73111	30.932	ng/u1	-0.01
15) 4-Methylphenol-d8	8.963	113	82049	31.234	ng/u1	-0.02
21) Nitrobenzene-d5	9.439	128	38507	31.323	ng/u1	-0.02
24) 2-Nitrophenol-d4	10.179	143	43264	32.193	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.719	165	83298	28.582	ng/u1	-0.01
31) 4-Chloroaniline-d4	11.231	131	105387	27.465	ng/u1	-0.01
46) Dimethylphthalate-d6	14.297	166	266145	28.217	ng/u1	-0.02
49) Acenaphthylene-d8	14.602	160	337344	28.798	ng/u1	-0.01
54) 4-Nitrophenol-d4	15.067	143	39656	25.854	ng/u1	-0.01
60) Fluorene-d10	15.895	176	236945	28.039	ng/u1	-0.01
65) 4,6-Dinitro-2-methylph...	15.995	200	35502	24.409	ng/u1	-0.02
73) Anthracene-d10	17.751	188	372888	28.001	ng/u1	-0.01
81) Pyrene-d10	20.030	212	435116	27.095	ng/u1	-0.01
92) Benzo(a)pyrene-d12	25.235	264	399569	28.236	ng/u1	-0.02

Target Compounds	Qvalue

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

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