

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011722\
 Data File : BG052028.D
 Acq On : 17 Jan 2022 17:30
 Operator : CG/JU
 Sample : N1169-14
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLY1

Quant Time: Jan 18 00:01:39 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG010622.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jan 06 14:43:02 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.249	152	23355	20.000	ng/ul	0.00
20) Naphthalene-d8	11.086	136	106329	20.000	ng/ul	0.00
38) Acenaphthene-d10	14.893	164	80101	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.642	188	190589	20.000	ng/ul	0.00
79) Chrysene-d12	21.966	240	170955	20.000	ng/ul	# 0.00
88) Perylene-d12	25.461	264	173206	20.000	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.567	96	3788	5.796	ng/uL	0.00
4) Pyridine-d5	4.002	84	47825	23.538	ng/ul	0.00
7) Phenol-d5	7.385	99	84559	36.993	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.550	67	52556	36.219	ng/ul	0.00
11) 2-Chlorophenol-d4	7.773	132	58154	38.426	ng/ul	0.00
15) 4-Methylphenol-d8	8.948	113	66281	37.301	ng/ul	0.00
21) Nitrobenzene-d5	9.418	128	32380	38.539	ng/ul	0.00
24) 2-Nitrophenol-d4	10.152	143	38135	41.040	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.699	165	69902	38.486	ng/ul	0.00
31) 4-Chloroaniline-d4	11.210	131	93705	37.046	ng/ul	0.00
46) Dimethylphthalate-d6	14.282	166	256479	38.737	ng/ul	0.00
49) Acenaphthylene-d8	14.588	160	319093	39.898	ng/ul	0.00
54) 4-Nitrophenol-d4	15.069	143	38361	34.654	ng/ul	0.00
60) Fluorene-d10	15.886	176	225947	39.141	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.991	200	37758	31.404	ng/ul	0.00
73) Anthracene-d10	17.742	188	378306	41.843	ng/ul	0.00
81) Pyrene-d10	20.021	212	422590	42.951	ng/ul	0.00
92) Benzo(a)pyrene-d12	25.220	264	377055	41.351	ng/ul	0.03
Target Compounds					Qvalue	
8) Phenol	7.409	94	3681	1.578	ng/ul#	75

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

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