

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011722\
 Data File : BG052045.D
 Acq On : 18 Jan 2022 9:51
 Operator : CG/JU
 Sample : N1169-17
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLY4

Quant Time: Jan 18 10:38:43 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	8.247	152	26722	20.000	ng/u1	0.00
20) Naphthalene-d8	11.090	136	115081	20.000	ng/u1	# 0.00
38) Acenaphthene-d10	14.891	164	84960	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.640	188	211830	20.000	ng/u1	# 0.00
79) Chrysene-d12	21.958	240	207322	20.000	ng/u1	# 0.00
88) Perylene-d12	25.447	264	210765	20.000	ng/u1	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.571	96	2365	3.162	ng/uL	0.00
4) Pyridine-d5	4.006	84	12117	5.212	ng/u1	0.00
7) Phenol-d5	7.390	99	42595	16.286	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.554	67	26247	15.809	ng/u1	0.00
11) 2-Chlorophenol-d4	7.771	132	29092	16.801	ng/u1	0.00
15) 4-Methylphenol-d8	8.946	113	31703	15.593	ng/u1	0.00
21) Nitrobenzene-d5	9.422	128	16425	18.062	ng/u1	0.00
24) 2-Nitrophenol-d4	10.151	143	18329	18.225	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.703	165	33163	16.870	ng/u1	0.00
31) 4-Chloroaniline-d4	11.214	131	42931	15.682	ng/u1	0.00
46) Dimethylphthalate-d6	14.280	166	121238	17.264	ng/u1	0.00
49) Acenaphthylene-d8	14.586	160	149970	17.679	ng/u1	0.00
54) 4-Nitrophenol-d4	15.073	143	15150	12.903	ng/u1	0.00
60) Fluorene-d10	15.878	176	109599	17.900	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.984	200	18225	13.638	ng/u1	0.00
73) Anthracene-d10	17.734	188	184735	18.384	ng/u1	0.00
81) Pyrene-d10	20.014	212	225787	18.923	ng/u1	0.00
92) Benzo(a)pyrene-d12	25.207	264	207993	18.746	ng/u1	0.01
Target Compounds						
					Qvalue	
8) Phenol	7.407	94	3280	1.229	ng/u1	94

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

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