

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG011722\
 Data File : BG052025.D
 Acq On : 17 Jan 2022 15:36
 Operator : CG/JU
 Sample : N1169-07
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 BGLX3

Quant Time: Jan 18 00:00:51 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	8.252	152	28370	20.000	ng/u1	0.00
20) Naphthalene-d8	11.089	136	125178	20.000	ng/u1	0.00
38) Acenaphthene-d10	14.896	164	76741	20.000	ng/u1	0.00
64) Phenanthrene-d10	17.645	188	159569	20.000	ng/u1	0.00
79) Chrysene-d12	21.969	240	162639	20.000	ng/u1	# 0.01
88) Perylene-d12	25.470	264	172449	20.000	ng/u1	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.570	96	1738	2.189	ng/uL	0.00
4) Pyridine-d5	0.000	84	0d	0.000	ng/u1	
7) Phenol-d5	7.388	99	48566	17.491	ng/u1	0.00
9) Bis-(2-Chloroethyl)eth...	7.553	67	35938	20.389	ng/u1	0.00
11) 2-Chlorophenol-d4	7.776	132	38024	20.683	ng/u1	0.00
15) 4-Methylphenol-d8	8.951	113	41358	19.161	ng/u1	0.00
21) Nitrobenzene-d5	9.421	128	21166	21.399	ng/u1	0.00
24) 2-Nitrophenol-d4	10.149	143	16868	15.419	ng/u1	0.00
28) 2,4-Dichlorophenol-d3	10.702	165	43319	20.259	ng/u1	0.00
31) 4-Chloroaniline-d4	11.224	131	358	0.120	ng/u1	0.01
46) Dimethylphthalate-d6	14.285	166	87577	13.806	ng/u1	0.00
49) Acenaphthylene-d8	14.590	160	172211	22.475	ng/u1	0.00
54) 4-Nitrophenol-d4	15.072	143	6900	6.506	ng/u1	0.00
60) Fluorene-d10	15.883	176	115843	20.946	ng/u1	0.00
65) 4,6-Dinitro-2-methylph...	15.989	200	7649	7.599	ng/u1	0.00
73) Anthracene-d10	17.745	188	180213	23.807	ng/u1	0.00
81) Pyrene-d10	20.024	212	209583	22.391	ng/u1	0.01
92) Benzo(a)pyrene-d12	25.223	264	201842	22.233	ng/u1	0.03

Target Compounds Qvalue

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

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