

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG071124\
 Data File : BG062014.D
 Acq On : 11 Jul 2024 17:44
 Operator : MA/JU
 Sample : PB161825BL
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SBLK825

Quant Time: Jul 12 00:07:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\SFAM-EPA-BG070224.MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Thu Jul 11 12:11:32 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.041	152	46774	20.000	ng/ul	0.00
20) Naphthalene-d8	10.850	136	220143	20.000	ng/ul	# 0.00
38) Acenaphthene-d10	14.674	164	174470	20.000	ng/ul	0.00
64) Phenanthrene-d10	17.412	188	460000	20.000	ng/ul	# 0.00
79) Chrysene-d12	21.672	240	429448	20.000	ng/ul	0.00
88) Perylene-d12	24.874	264	489165	20.000	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.441	96	6531	5.247	ng/uL	0.00
4) Pyridine-d5	3.858	84	91775	23.980	ng/ul	0.00
7) Phenol-d5	7.195	99	130044	25.405	ng/ul	0.00
9) Bis-(2-Chloroethyl)eth...	7.365	67	77120	23.791	ng/ul	0.00
11) 2-Chlorophenol-d4	7.571	132	90413	25.750	ng/ul	0.00
15) 4-Methylphenol-d8	8.740	113	98736	24.499	ng/ul	0.00
21) Nitrobenzene-d5	9.210	128	50093	29.919	ng/ul	0.00
24) 2-Nitrophenol-d4	9.933	143	57066	29.847	ng/ul	0.00
28) 2,4-Dichlorophenol-d3	10.468	165	101790	27.601	ng/ul	0.00
31) 4-Chloroaniline-d4	10.985	131	120069	22.425	ng/ul	0.00
46) Dimethylphthalate-d6	14.069	166	385914	26.905	ng/ul	0.00
49) Acenaphthylene-d8	14.369	160	405899	27.063	ng/ul	0.00
54) 4-Nitrophenol-d4	14.868	143	59929	26.163	ng/ul	0.00
60) Fluorene-d10	15.662	176	341316	28.266	ng/ul	0.00
65) 4,6-Dinitro-2-methylph...	15.773	200	54096	22.392	ng/ul	0.00
73) Anthracene-d10	17.512	188	580902	27.048	ng/ul	0.00
81) Pyrene-d10	19.798	212	700508	27.732	ng/ul	0.00
92) Benzo(a)pyrene-d12	24.651	264	713393	29.406	ng/ul	-0.01

Target Compounds	Qvalue

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

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