

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040624\
 Data File : BG060818.D
 Acq On : 6 Apr 2024 19:37
 Operator : MA/JU
 Sample : P2000-05 20X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 60153

Quant Time: Apr 08 02:10:26 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG032924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Mar 30 03:12:26 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.953	152	79426	20.000	ng	0.00
21) Naphthalene-d8	10.767	136	305987	20.000	ng	0.00
39) Acenaphthene-d10	14.592	164	212033	20.000	ng	0.00
64) Phenanthrene-d10	17.336	188	453924	20.000	ng	0.00
76) Chrysene-d12	21.596	240	436325	20.000	ng	0.00
86) Perylene-d12	24.733	264	528243	20.000	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.544	112	11202	2.350	ng	0.00
7) Phenol-d6	7.131	99	12066	1.705	ng	0.02
23) Nitrobenzene-d5	9.128	82	20518	3.826	ng	0.02
42) 2,4,6-Tribromophenol	16.085	330	14814	6.155	ng	0.00
45) 2-Fluorobiphenyl	13.218	172	47639	3.297	ng	0.00
79) Terphenyl-d14	19.951	244	75017	3.027	ng	-0.01
Target Compounds						
						Qvalue
10) Phenol	7.172	94	2598606	359.783	ng	90
15) Benzyl Alcohol	8.265	79	61400	10.710	ng	99
17) 2-Methylphenol	8.411	107	1023036	187.625	ng	97
20) 3+4-Methylphenols	8.817	107	10902254	1460.143	ng	# 73
27) 2,4-Dimethylphenol	9.974	122	4306234	867.945	ng	100
31) Naphthalene	10.814	128	43120	2.605	ng	# 84

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

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