

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040624\
 Data File : BG060817.D
 Acq On : 6 Apr 2024 18:56
 Operator : MA/JU
 Sample : P2000-04
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 60152

Quant Time: Apr 08 02:09:50 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	7.955	152	75472	20.000	ng	0.00
21) Naphthalene-d8	10.752	136	295023	20.000	ng	0.00
39) Acenaphthene-d10	14.583	164	212214	20.000	ng	0.00
64) Phenanthrene-d10	17.332	188	434031	20.000	ng	0.00
76) Chrysene-d12	21.598	240	404335	20.000	ng	# 0.00
86) Perylene-d12	24.741	264	498271	20.000	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.546	112	263419	58.144	ng	0.01
7) Phenol-d6	7.121	99	213388	31.731	ng	0.00
23) Nitrobenzene-d5	9.112	82	539459	104.339	ng	0.00
42) 2,4,6-Tribromophenol	16.081	330	377201	156.582	ng	0.00
45) 2-Fluorobiphenyl	13.214	172	1228744	84.977	ng	0.00
79) Terphenyl-d14	19.959	244	1778146	77.436	ng	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.472	88	14303	7.700	ng	# 76
3) Pyridine	3.860	79	61991	11.064	ng	93
10) Phenol	7.150	94	84547	12.319	ng	95
17) 2-Methylphenol	8.396	107	56553	10.915	ng	98
20) 3+4-Methylphenols	8.725	107	1033743	145.703	ng	97
22) Acetophenone	8.778	105	715909	88.005	ng	# 97
27) 2,4-Dimethylphenol	9.929	122	126724	26.491	ng	98
31) Naphthalene	10.799	128	244693	15.333	ng	97
50) Dimethylphthalate	14.048	163	35916	2.141	ng	# 94
66) n-Nitrosodiphenylamine	15.846	169	45052	3.632	ng	94
74) Di-n-butylphthalate	18.319	149	1152996	45.756	ng	# 97

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

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