

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062916\
 Data File : BG022713.D
 Acq On : 30 Jun 2016 3:58
 Operator : UM/SJ
 Sample : H3828-09
 Misc :
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S8-0579

Quant Time: Jun 30 14:01:50 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 20:14:59 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	118435	20.00	ng	-0.02
21) Naphthalene-d8	11.00	136	560538	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	554649	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	958122	20.00	ng	0.00
75) Chrysene-d12	21.85	240	1045648	20.00	ng	0.00
86) Perylene-d12	25.22	264	1080568	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	458318	62.07	ng	0.00
7) Phenol-d6	7.33	99	541477	52.03	ng	0.03
23) Nitrobenzene-d5	9.34	82	1091351	82.40	ng	0.00
41) 2,4,6-Tribromophenol	16.30	330	880784	100.96	ng	0.01
44) 2-Fluorobiphenyl	13.43	172	2213754	63.30	ng	0.00
78) Terphenyl-d14	20.16	244	2658501	67.36	ng	0.00
Target Compounds						Qvalue
9) Benzaldehyde	7.30	77	1316620	359.46	ng	# 88
15) Benzyl Alcohol	8.41	79	186404	19.36	ng	94
17) 2-Methylphenol	8.63	107	31444	4.34	ng	# 86
20) 3+4-Methylphenols	8.95	107	60487	6.06	ng	97
22) Acetophenone	9.02	105	5802908	358.10	ng	# 67
27) 2,4-Dimethylphenol	10.17	122	310489	30.82	ng	91
31) Naphthalene	11.04	128	1260080	44.72	ng	98
37) 2-Methylnaphthalene	12.64	142	282340	12.92	ng	98

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

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