

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062916\
 Data File : BG022712.D
 Acq On : 30 Jun 2016 3:19
 Operator : UM/SJ
 Sample : H3828-08
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 S8-0578

Quant Time: Jun 30 13:55:32 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	128833	20.00	ng	0.00
21) Naphthalene-d8	10.99	136	602880	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	565687	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	1037883	20.00	ng	0.00
75) Chrysene-d12	21.85	240	1101439	20.00	ng	0.00
86) Perylene-d12	25.22	264	1140691	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	507272	63.15	ng	0.00
7) Phenol-d6	7.33	99	527243	46.57	ng	0.03
23) Nitrobenzene-d5	9.35	82	1169034	82.07	ng	0.00
41) 2,4,6-Tribromophenol	16.30	330	931553	104.70	ng	0.01
44) 2-Fluorobiphenyl	13.43	172	2313548	64.86	ng	0.00
78) Terphenyl-d14	20.16	244	2616426	62.94	ng	0.00
Target Compounds						Qvalue
15) Benzyl Alcohol	8.41	79	164968	15.75	ng	98
20) 3+4-Methylphenols	8.96	107	50320	4.63	ng	# 74
22) Acetophenone	9.02	105	5534714	317.56	ng	# 70
27) 2,4-Dimethylphenol	10.17	122	298504	27.55	ng	# 93
31) Naphthalene	11.04	128	1227180	40.49	ng	96
32) Benzoic acid	10.31	122	58223	12.46	ng	93
37) 2-Methylnaphthalene	12.64	142	281982	12.00	ng	99

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

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