

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG060917\
 Data File : BG027359.D
 Acq On : 9 Jun 2017 22:45
 Operator : SJ/MA
 Sample : I3561-01
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PURGE-WATER-NRL-65

Quant Time: Jun 10 06:46:46 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 16:00:35 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.55	152	46035	20.00	ng	-0.02
21) Naphthalene-d8	11.37	136	221253	20.00	ng	-0.02
38) Acenaphthene-d10	15.14	164	134858	20.00	ng	-0.02
63) Phenanthrene-d10	17.88	188	314110	20.00	ng	-0.03
75) Chrysene-d12	22.27	240	369947	20.00	ng	-0.04
86) Perylene-d12	25.95	264	360515	20.00	ng	-0.07
System Monitoring Compounds						
5) 2-Fluorophenol	6.14	112	159326	52.81	ng	-0.02
7) Phenol-d6	7.69	99	144407	32.61	ng	-0.01
23) Nitrobenzene-d5	9.72	82	306125	87.01	ng	-0.02
41) 2,4,6-Tribromophenol	16.62	330	203490	114.59	ng	-0.03
44) 2-Fluorobiphenyl	13.76	172	659915	68.46	ng	-0.03
78) Terphenyl-d14	20.46	244	259261	17.88	ng	-0.04
Target Compounds						
9) Benzaldehyde	7.70	77	13498	4.409	ng	Qvalue 97
10) Phenol	7.71	94	29766	6.573	ng	81
15) Benzyl Alcohol	8.78	79	431424	138.271	ng	# 86
22) Acetophenone	9.38	105	11486	2.030	ng	# 84
32) Benzoic acid	10.59	122	12889	12.387	ng	92
49) Dimethylphthalate	14.57	163	124962	11.681	ng	97

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

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