

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122117\
 Data File : BG031708.D
 Acq On : 22 Dec 2017 13:08
 Operator : SJ/JU
 Sample : I6971-12
 Misc :
 ALS Vial : 33 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 121517-DBRI-E-D-S

Quant Time: Dec 22 14:04:33 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122117.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 22 11:08:55 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.83	152	55359	20.00	ng	-0.01
21) Naphthalene-d8	11.71	136	230445	20.00	ng	-0.01
38) Acenaphthene-d10	15.46	164	164433	20.00	ng	-0.01
63) Phenanthrene-d10	18.19	188	431342	20.00	ng	-0.01
75) Chrysene-d12	22.55	240	499891	20.00	ng	-0.02
86) Perylene-d12	26.34	264	523997	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	6.24	112	122145	40.19	ng	0.00
7) Phenol-d6	7.92	99	93310	23.65	ng	-0.01
23) Nitrobenzene-d5	10.03	82	282303	92.50	ng	-0.01
41) 2,4,6-Tribromophenol	16.93	330	243786	97.16	ng	-0.01
44) 2-Fluorobiphenyl	14.09	172	1070598	81.72	ng	-0.01
78) Terphenyl-d14	20.73	244	1875680	85.72	ng	-0.01
Target Compounds						
37) 2-Methylnaphthalene	13.32	142	25307	2.684	ng	Qvalue # 93
51) Acenaphthene	15.52	154	30691	2.734	ng	98
57) Fluorene	16.50	166	30504	2.182	ng	99
70) Phenanthrene	18.24	178	131733	5.397	ng	99

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

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