

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG122217\
 Data File : BG031726.D
 Acq On : 23 Dec 2017 00:25
 Operator : SJ/JU
 Sample : I7004-05
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PR-MW-1-121817

Quant Time: Dec 26 13:41:32 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	51698	20.00	ng	-0.01
21) Naphthalene-d8	11.72	136	210862	20.00	ng	0.00
38) Acenaphthene-d10	15.46	164	144607	20.00	ng	0.00
63) Phenanthrene-d10	18.20	188	357504	20.00	ng	0.00
75) Chrysene-d12	22.55	240	427964	20.00	ng	-0.02
86) Perylene-d12	26.34	264	463177	20.00	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	6.25	112	180771	63.69	ng	0.00
7) Phenol-d6	7.92	99	146938	39.87	ng	0.00
23) Nitrobenzene-d5	10.03	82	299180	107.14	ng	0.00
41) 2,4,6-Tribromophenol	16.93	330	358114	162.30	ng	0.00
44) 2-Fluorobiphenyl	14.09	172	1076813	93.46	ng	0.00
78) Terphenyl-d14	20.73	244	1849685	98.74	ng	0.00
Target Compounds						
10) Phenol	7.94	94	8041	2.161	ng	Qvalue # 5
20) 3+4-Methylphenols	9.58	107	8542	2.258	ng	96
27) 2,4-Dimethylphenol	10.86	122	7536	2.373	ng	96
31) Naphthalene	11.77	128	1030501	96.836	ng	98
37) 2-Methylnaphthalene	13.33	142	543261	62.973	ng	91
45) 1,1'-Biphenyl	14.30	154	58223	4.714	ng	93

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

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