

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101617\
 Data File : BG029292.D
 Acq On : 17 Oct 2017 4:43
 Operator : SJ/JU
 Sample : I5906-13
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 WH-TWS-03

Quant Time: Oct 17 05:37:23 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG101617.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Oct 16 17:32:15 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	71404	20.00	ng	0.00
21) Naphthalene-d8	10.99	136	340259	20.00	ng	0.00
38) Acenaphthene-d10	14.79	164	209850	20.00	ng	0.00
63) Phenanthrene-d10	17.53	188	464650	20.00	ng	0.00
75) Chrysene-d12	21.79	240	508922	20.00	ng	0.00
86) Perylene-d12	25.11	264	521559	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.72	112	559486	130.90	ng	0.00
7) Phenol-d6	7.33	99	692251	115.38	ng	0.00
23) Nitrobenzene-d5	9.34	82	473731	88.47	ng	0.00
41) 2,4,6-Tribromophenol	16.27	330	323008	119.22	ng	0.00
44) 2-Fluorobiphenyl	13.41	172	1173450	82.01	ng	0.00
78) Terphenyl-d14	20.12	244	1737862	77.68	ng	0.00

Target Compounds

						Qvalue
3) Pyridine	4.01	79	12187	2.413	ng	94
10) Phenol	7.36	94	1005714	155.487	ng	95
17) 2-Methylphenol	8.61	107	398536	92.753	ng	97
20) 3+4-Methylphenols	8.96	107	1326901	219.169	ng	88
27) 2,4-Dimethylphenol	10.17	122	582105	111.815	ng	93
31) Naphthalene	11.05	128	4635447	273.351	ng	# 84
37) 2-Methylnaphthalene	12.63	142	449153	36.342	ng	94
45) 1,1'-Biphenyl	13.63	154	97079	5.382	ng	92
48) Acenaphthylene	14.51	152	196850	9.560	ng	98
51) Acenaphthene	14.85	154	194621	14.527	ng	96
54) Dibenzofuran	15.18	168	330523	17.282	ng	100
57) Fluorene	15.83	166	178650	11.307	ng	93
70) Phenanthrene	17.57	178	448664	18.691	ng	99
71) Anthracene	17.66	178	77611	3.220	ng	95
72) Carbazole	17.93	167	909788	39.293	ng	98
74) Fluoranthene	19.56	202	101599	3.619	ng	98
77) Pyrene	19.92	202	62892	2.037	ng	98

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

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