

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041416\
 Data File : BG021657.D
 Acq On : 14 Apr 2016 17:55
 Operator : UM/SJ
 Sample : H1739-06DL 5X
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Apr 15 01:55:22 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG041316.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 13 15:48:13 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.22	152	31637	20.00	ng	0.00
21) Naphthalene-d8	11.03	136	146240	20.00	ng	0.00
38) Acenaphthene-d10	14.82	164	96874	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	238724	20.00	ng	0.00
75) Chrysene-d12	21.83	240	269804	20.00	ng	0.00
86) Perylene-d12	25.16	264	260974	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.76	112	50243	26.45	ng	0.00
7) Phenol-d6	7.36	99	75093	26.46	ng	0.00
23) Nitrobenzene-d5	9.38	82	46615	16.38	ng	0.00
41) 2,4,6-Tribromophenol	16.30	330	28968	27.75	ng	0.00
44) 2-Fluorobiphenyl	13.45	172	115320	17.44	ng	0.00
78) Terphenyl-d14	20.14	244	188831	17.46	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	7.78	128	59880	26.29	ng	93
17) 2-Methylphenol	8.65	107	52843	25.04	ng	90
20) 3+4-Methylphenols	8.97	107	86481	29.95	ng	93
26) 2-Nitrophenol	10.13	139	21265	18.28	ng	96
27) 2,4-Dimethylphenol	10.19	122	45544	17.24	ng	93
29) 2,4-Dichlorophenol	10.67	162	49385	21.93	ng	97
36) 4-Chloro-3-methylphenol	12.27	107	41401	14.62	ng	95
42) 2,4,6-Trichlorophenol	13.26	196	52046	26.96	ng	97
43) 2,4,5-Trichlorophenol	13.33	196	34697	16.66	ng	95
53) 2,4-Dinitrophenol	14.92	184	10823	23.43	ng	93
55) 4-Nitrophenol	15.01	139	39473	26.03	ng	94
64) 4,6-Dinitro-2-methylphenol	15.92	198	8232	12.75	ng	94
69) Pentachlorophenol	17.20	266	37070	24.05	ng	93

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

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