

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG041416\
 Data File : BG021656.D
 Acq On : 14 Apr 2016 17:16
 Operator : UM/SJ
 Sample : H1739-06
 Misc :
 ALS Vial : 43 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Apr 15 01:47:05 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	30161	20.00	ng	0.00
21) Naphthalene-d8	11.03	136	132232	20.00	ng	0.00
38) Acenaphthene-d10	14.82	164	88431	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	218904	20.00	ng	0.00
75) Chrysene-d12	21.83	240	251457	20.00	ng	0.00
86) Perylene-d12	25.16	264	246161	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.76	112	253133	139.76	ng	0.00
7) Phenol-d6	7.36	99	377377	139.49	ng	0.00
23) Nitrobenzene-d5	9.38	82	229648	89.27	ng	0.00
41) 2,4,6-Tribromophenol	16.30	330	153181	160.73	ng	0.00
44) 2-Fluorobiphenyl	13.45	172	529276	87.69	ng	0.00
78) Terphenyl-d14	20.15	244	893258	88.64	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	7.78	128	293836	135.33	ng	97
17) 2-Methylphenol	8.65	107	256589	127.54	ng	94
20) 3+4-Methylphenols	8.97	107	418415	151.99	ng	95
26) 2-Nitrophenol	10.14	139	116952	111.21	ng	98
27) 2,4-Dimethylphenol	10.19	122	222411	93.10	ng	98
29) 2,4-Dichlorophenol	10.67	162	240600	118.13	ng	97
36) 4-Chloro-3-methylphenol	12.28	107	204707	79.93	ng	99
42) 2,4,6-Trichlorophenol	13.26	196	260920	148.06	ng	97
43) 2,4,5-Trichlorophenol	13.33	196	165615	87.09	ng	98
53) 2,4-Dinitrophenol	14.92	184	76938	131.85	ng	97
55) 4-Nitrophenol	15.02	139	220008	158.91	ng	93
64) 4,6-Dinitro-2-methylphenol	15.93	198	53929	57.90	ng	97
69) Pentachlorophenol	17.20	266	221410	156.64	ng	94

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

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