

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032017\
 Data File : BG026332.D
 Acq On : 20 Mar 2017 18:27
 Operator : SJ/MA
 Sample : I2194-06DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Mar 21 04:50:04 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG032017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 20 17:43:27 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.05	152	119340	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	493901	20.00	ng	0.00
38) Acenaphthene-d10	14.66	164	276426	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	660888	20.00	ng	0.00
75) Chrysene-d12	21.63	240	837247	20.00	ng	0.00
86) Perylene-d12	24.81	264	825167	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	204645	30.44	ng	0.00
7) Phenol-d6	7.21	99	276892	30.67	ng	0.00
23) Nitrobenzene-d5	9.20	82	149395	18.19	ng	0.00
41) 2,4,6-Tribromophenol	16.13	330	99158	31.32	ng	0.00
44) 2-Fluorobiphenyl	13.28	172	412327	20.92	ng	0.00
78) Terphenyl-d14	19.98	244	634765	18.41	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	7.61	128	259170	34.23	ng	83
10) Phenol	7.23	94	271705	28.21	ng	# 70
17) 2-Methylphenol	8.49	107	211266	30.88	ng	# 81
20) 3+4-Methylphenols	8.80	107	192211	20.41	ng	# 75
26) 2-Nitrophenol	9.96	139	157470	37.31	ng	# 64
27) 2,4-Dimethylphenol	10.01	122	196062	28.30	ng	86
29) 2,4-Dichlorophenol	10.50	162	198707	28.36	ng	95
32) Benzoic acid	10.11	122	88579	16.62	ng	# 74
36) 4-Chloro-3-methylphenol	12.12	107	55025	7.34	ng	# 79
42) 2,4,6-Trichlorophenol	13.09	196	183140	33.63	ng	95
43) 2,4,5-Trichlorophenol	13.16	196	237193	39.41	ng	94
53) 2,4-Dinitrophenol	14.76	184	85386	30.23	ng	# 81
55) 4-Nitrophenol	14.86	139	148792	33.89	ng	# 42
64) 4,6-Dinitro-2-methylphenol	15.77	198	113902	27.34	ng	83

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

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