

Data Path : Z:\HPCHEM1\BNA G\DATA\BG032017\
 Data File : BG026331.D
 Acq On : 20 Mar 2017 17:47
 Operator : SJ/MA
 Sample : I2194-06
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Mar 21 04:48:02 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	129651	20.00	ng	0.00
21) Naphthalene-d8	10.85	136	538689	20.00	ng	0.00
38) Acenaphthene-d10	14.66	164	309483	20.00	ng	0.00
63) Phenanthrene-d10	17.39	188	729584	20.00	ng	0.00
75) Chrysene-d12	21.63	240	856619	20.00	ng	0.00
86) Perylene-d12	24.81	264	811704	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.62	112	1044738	143.02	ng	0.00
7) Phenol-d6	7.21	99	1414244	144.21	ng	0.00
23) Nitrobenzene-d5	9.20	82	793874	88.61	ng	0.00
41) 2,4,6-Tribromophenol	16.14	330	515717	145.51	ng	0.00
44) 2-Fluorobiphenyl	13.28	172	1944681	88.12	ng	0.00
78) Terphenyl-d14	19.99	244	2704642	76.68	ng	0.00
Target Compounds						
8) 2-Chlorophenol	7.62	128	1282674	155.94	ng	Qvalue 86
10) Phenol	7.24	94	1344747	128.50	ng	# 73
17) 2-Methylphenol	8.49	107	1076824	144.89	ng	# 82
20) 3+4-Methylphenols	8.81	107	1004512	98.17	ng	# 79
26) 2-Nitrophenol	9.96	139	830901	180.52	ng	# 67
27) 2,4-Dimethylphenol	10.02	122	987159	130.66	ng	85
29) 2,4-Dichlorophenol	10.50	162	1021780	133.73	ng	95
32) Benzoic acid	10.20	122	538742	92.67	ng	# 72
36) 4-Chloro-3-methylphenol	12.11	107	309819	37.91	ng	# 80
42) 2,4,6-Trichlorophenol	13.10	196	953484	156.39	ng	96
43) 2,4,5-Trichlorophenol	13.17	196	1193063	177.04	ng	# 92
53) 2,4-Dinitrophenol	14.77	184	546753	172.88	ng	# 80
55) 4-Nitrophenol	14.87	139	829360	168.72	ng	# 42
64) 4,6-Dinitro-2-methylphenol	15.78	198	647293	140.72	ng	85

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

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