

Data Path : Z:\HPCHEM1\BNA G\DATA\BG101617\
 Data File : BG029278.D
 Acq On : 16 Oct 2017 18:03
 Operator : SJ/JU
 Sample : I5266-06
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PT-ACIDS-WP

Quant Time: Oct 16 19:14:36 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	8.18	152	54350	20.00	ng	0.01
21) Naphthalene-d8	10.99	136	237543	20.00	ng	0.00
38) Acenaphthene-d10	14.79	164	151697	20.00	ng	0.00
63) Phenanthrene-d10	17.53	188	358687	20.00	ng	0.00
75) Chrysene-d12	21.79	240	445860	20.00	ng	0.00
86) Perylene-d12	25.11	264	456027	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.74	112	464523	142.78	ng	0.02
7) Phenol-d6	7.34	99	594150	130.11	ng	0.01
23) Nitrobenzene-d5	9.34	82	327706	87.67	ng	0.00
41) 2,4,6-Tribromophenol	16.27	330	253264	129.31	ng	0.00
44) 2-Fluorobiphenyl	13.42	172	854996	82.66	ng	0.00
78) Terphenyl-d14	20.11	244	1403017	71.59	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	7.75	128	675180	181.053	ng	94
10) Phenol	7.36	94	820715	166.700	ng	95
17) 2-Methylphenol	8.61	107	269867	82.515	ng	98
20) 3+4-Methylphenols	8.93	107	409436	88.848	ng	95
26) 2-Nitrophenol	10.10	139	187918	93.549	ng	89
27) 2,4-Dimethylphenol	10.15	122	327987	90.245	ng	94
29) 2,4-Dichlorophenol	10.64	162	584239	150.746	ng	94
36) 4-Chloro-3-methylphenol	12.25	107	195640	45.897	ng	# 87
42) 2,4,6-Trichlorophenol	13.23	196	300682	86.482	ng	99
43) 2,4,5-Trichlorophenol	13.30	196	233589	65.420	ng	96
55) 4-Nitrophenol	15.00	139	397431	180.056	ng	# 82
64) 4,6-Dinitro-2-methylphenol	15.91	198	287895	131.442	ng	87
69) Pentachlorophenol	17.18	266	345566	129.030	ng	97

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

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