

Data Path : Z:\HPCHEM1\BNA G\DATA\BG031918\
 Data File : BG033243.D
 Acq On : 19 Mar 2018 17:08
 Operator : SJ/JU
 Sample : J1888-06DL 5X
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PT-ACIDS-WPDL

Quant Time: Mar 19 17:46:19 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG031918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 19 16:48:53 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.90	152	31644	20.00	ng	0.00
21) Naphthalene-d8	11.78	136	132121	20.00	ng	0.00
38) Acenaphthene-d10	15.51	164	106506	20.00	ng	0.00
63) Phenanthrene-d10	18.24	188	295382	20.00	ng	0.00
75) Chrysene-d12	22.59	240	309852	20.00	ng	0.00
86) Perylene-d12	26.40	264	318456	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	6.32	112	43817	25.96	ng	0.00
7) Phenol-d6	8.01	99	76119	26.25	ng	0.00
23) Nitrobenzene-d5	10.10	82	51181	17.80	ng	0.00
41) 2,4,6-Tribromophenol	16.98	330	42759	26.84	ng	0.00
44) 2-Fluorobiphenyl	14.14	172	137146	18.59	ng	0.00
78) Terphenyl-d14	20.76	244	310776	21.45	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	8.45	128	54327	28.160	ng	92
10) Phenol	8.04	94	60919	21.279	ng	91
17) 2-Methylphenol	9.35	107	16824	8.627	ng	96
20) 3+4-Methylphenols	9.66	107	40213	14.482	ng	90
26) 2-Nitrophenol	10.88	139	21499	18.449	ng	# 90
27) 2,4-Dimethylphenol	10.90	122	47811	28.242	ng	83
29) 2,4-Dichlorophenol	11.41	162	63144	30.330	ng	94
36) 4-Chloro-3-methylphenol	12.99	107	79178	29.159	ng	90
42) 2,4,6-Trichlorophenol	13.96	196	68008	30.341	ng	95
43) 2,4,5-Trichlorophenol	14.04	196	21635	8.166	ng	87
55) 4-Nitrophenol	15.69	139	32552	23.298	ng	97
64) 4,6-Dinitro-2-methylphenol	16.60	198	23305	13.189	ng	97
69) Pentachlorophenol	17.88	266	26731	10.638	ng	94

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

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