

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080216\
 Data File : BG023395.D
 Acq On : 2 Aug 2016 10:41
 Operator : UM/SJ
 Sample : H4145-12
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 9535

Quant Time: Aug 02 12:57:56 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG080116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 01 19:22:14 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.13	152	130848	20.00	ng	0.01
21) Naphthalene-d8	10.96	136	637988	20.00	ng	0.01
38) Acenaphthene-d10	14.79	164	472423	20.00	ng	0.00
63) Phenanthrene-d10	17.55	188	1141164	20.00	ng	0.00
75) Chrysene-d12	21.86	240	1044367	20.00	ng	0.00
86) Perylene-d12	25.25	264	1049858	20.00	ng	0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.67	112	1169463	150.44	ng	0.01
7) Phenol-d6	7.29	99	1816295	150.01	ng	0.02
23) Nitrobenzene-d5	9.31	82	1154632	89.97	ng	0.01
41) 2,4,6-Tribromophenol	16.29	330	954799	138.18	ng	0.00
44) 2-Fluorobiphenyl	13.41	172	2372828	84.72	ng	0.01
78) Terphenyl-d14	20.16	244	3196276	98.91	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	7.69	128	473887	53.06	ng	98
10) Phenol	7.32	94	1804296	139.30	ng	85
20) 3+4-Methylphenols	8.91	107	1856982	151.70	ng	85
26) 2-Nitrophenol	10.07	139	509067	84.88	ng	# 80
29) 2,4-Dichlorophenol	10.61	162	1267340	123.42	ng	97
36) 4-Chloro-3-methylphenol	12.24	107	959317	71.11	ng	90
42) 2,4,6-Trichlorophenol	13.23	196	1595764	156.37	ng	97
43) 2,4,5-Trichlorophenol	13.30	196	446876	42.53	ng	# 95
53) 2,4-Dinitrophenol	14.91	184	436003	79.24	ng	87
55) 4-Nitrophenol	15.02	139	1075610	141.49	ng	# 76
64) 4,6-Dinitro-2-methylphenol	15.92	198	526302	67.83	ng	96
69) Pentachlorophenol	17.20	266	1106900	130.48	ng	96

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

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