

Data Path : Z:\HPCHEM1\BNA G\DATA\BG080216\
 Data File : BG023397.D
 Acq On : 2 Aug 2016 12:01
 Operator : UM/SJ
 Sample : H4145-12DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 9535DL

Quant Time: Aug 02 15:56:13 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	8.12	152	113803	20.00	ng	0.00
21) Naphthalene-d8	10.95	136	508981	20.00	ng	0.00
38) Acenaphthene-d10	14.78	164	352469	20.00	ng	0.00
63) Phenanthrene-d10	17.54	188	900367	20.00	ng	0.00
75) Chrysene-d12	21.86	240	987328	20.00	ng	0.00
86) Perylene-d12	25.24	264	995679	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.66	112	199707	29.54	ng	0.00
7) Phenol-d6	7.27	99	309704	29.41	ng	0.00
23) Nitrobenzene-d5	9.29	82	193424	18.89	ng	0.00
41) 2,4,6-Tribromophenol	16.28	330	152717	29.62	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	462559	22.13	ng	0.00
78) Terphenyl-d14	20.15	244	844965	9.86	ng	0.00
Target Compounds						
8) 2-Chlorophenol	7.68	128	79416	10.22	ng	Qvalue 90
10) Phenol	7.30	94	319816	28.39	ng	# 79
20) 3+4-Methylphenols	8.89	107	313306	29.43	ng	# 78
26) 2-Nitrophenol	10.06	139	73614	15.38	ng	# 78
29) 2,4-Dichlorophenol	10.60	162	202202	24.68	ng	97
36) 4-Chloro-3-methylphenol	12.24	107	147194	13.68	ng	# 94
42) 2,4,6-Trichlorophenol	13.21	196	280644	36.86	ng	97
43) 2,4,5-Trichlorophenol	13.30	196	66630	8.50	ng	94
53) 2,4-Dinitrophenol	14.90	184	47329	11.53	ng	96
55) 4-Nitrophenol	15.02	139	155886	27.48	ng	# 75
64) 4,6-Dinitro-2-methylphenol	15.91	198	69197	11.30	ng	96
69) Pentachlorophenol	17.19	266	157965	23.60	ng	97

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

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