

Data Path : Z:\HPCHEM1\BNA F\DATA\BF013017\
 Data File : BF092634.D
 Acq On : 30 Jan 2017 15:27
 Operator : SJ/MA
 Sample : I1305-22DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 0035DL

Quant Time: Jan 30 15:55:05 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF013017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 30 14:50:05 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.96	152	162277	20.00	ng	-0.01
21) Naphthalene-d8	8.25	136	647069	20.00	ng	-0.01
38) Acenaphthene-d10	10.01	164	364275	20.00	ng	0.00
63) Phenanthrene-d10	11.49	188	596946	20.00	ng	0.00
75) Chrysene-d12	14.13	240	559817	20.00	ng	-0.01
86) Perylene-d12	15.61	264	461561	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.56	112	305624	32.31	ng	-0.01
7) Phenol-d6	6.59	99	378179	31.07	ng	-0.01
23) Nitrobenzene-d5	7.53	82	246784	21.24	ng	-0.01
41) 2,4,6-Tribromophenol	10.80	330	87024	33.03	ng	-0.01
44) 2-Fluorobiphenyl	9.33	172	480615	23.90	ng	0.00
78) Terphenyl-d14	13.08	244	488759	20.51	ng	0.00
Target Compounds						Qvalue
8) 2-Chlorophenol	6.75	128	185137	17.74	ng	86
17) 2-Methylphenol	7.22	107	154973	17.31	ng	95
20) 3+4-Methylphenols	7.38	107	449014	41.95	ng	# 61
27) 2,4-Dimethylphenol	7.90	122	138167	13.87	ng	95
29) 2,4-Dichlorophenol	8.11	162	300653	32.36	ng	91
36) 4-Chloro-3-methylphenol	8.81	107	194888	20.12	ng	93
42) 2,4,6-Trichlorophenol	9.24	196	174727	26.00	ng	92
43) 2,4,5-Trichlorophenol	9.29	196	147159	19.99	ng	# 84
53) 2,4-Dinitrophenol	10.06	184	64489	27.24	ng	# 66
55) 4-Nitrophenol	10.13	139	89758	20.55	ng	# 66
64) 4,6-Dinitro-2-methylphenol	10.60	198	61433	18.64	ng	94
69) Pentachlorophenol	11.30	266	41328	19.02	ng	99

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

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