

Data Path : Z:\HPCHEM1\BNA F\DATA\BF012417\
 Data File : BF092427.D
 Acq On : 24 Jan 2017 14:23
 Operator : UM/SJ
 Sample : I1305-22DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 0035DL

Quant Time: Jan 24 14:57:49 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF012417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 24 13:48:07 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.97	152	254088	20.00	ng	0.00
21) Naphthalene-d8	8.26	136	967995	20.00	ng	-0.01
38) Acenaphthene-d10	10.03	164	572765	20.00	ng	0.00
63) Phenanthrene-d10	11.52	188	866648	20.00	ng	-0.01
75) Chrysene-d12	14.16	240	818889	20.00	ng	-0.01
86) Perylene-d12	15.64	264	706204	20.00	ng	0.00
System Monitoring Compounds						
5) 2-Fluorophenol	5.55	112	346746	21.23	ng	0.00
7) Phenol-d6	6.58	99	471999	22.69	ng	-0.01
23) Nitrobenzene-d5	7.53	82	255987	14.45	ng	-0.01
41) 2,4,6-Tribromophenol	10.82	330	87302	22.37	ng	0.00
44) 2-Fluorobiphenyl	9.34	172	513655	16.58	ng	-0.01
78) Terphenyl-d14	13.11	244	465390	15.13	ng	-0.01
Target Compounds						Qvalue
8) 2-Chlorophenol	6.74	128	188678	11.00	ng	86
17) 2-Methylphenol	7.21	107	159523	10.88	ng	97
20) 3+4-Methylphenols	7.37	107	503030	28.29	ng	# 63
27) 2,4-Dimethylphenol	7.90	122	153616	9.49	ng	97
29) 2,4-Dichlorophenol	8.11	162	317209	22.61	ng	95
36) 4-Chloro-3-methylphenol	8.80	107	179880	11.89	ng	91
42) 2,4,6-Trichlorophenol	9.25	196	202860	18.33	ng	90
43) 2,4,5-Trichlorophenol	9.29	196	132563	13.09	ng	# 89
53) 2,4-Dinitrophenol	10.08	184	58573	22.89	ng	# 76
55) 4-Nitrophenol	10.12	139	93499	12.63	ng	# 66
64) 4,6-Dinitro-2-methylphenol	10.61	198	59386	16.36	ng	92
69) Pentachlorophenol	11.32	266	77009	13.62	ng	97

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

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