

Data Path : Z:\HPCHEM1\BNA F\DATA\BF020916\
 Data File : BF084981.D
 Acq On : 10 Feb 2016 4:46
 Operator : SJ/IZ
 Sample : H1360-01
 Misc :
 ALS Vial : 42 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SW007

Quant Time: Feb 10 06:24:08 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF020116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 03 14:16:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.64	152	156726	20.00	ng	-0.05
21) Naphthalene-d8	7.92	136	568471	20.00	ng	-0.06
38) Acenaphthene-d10	9.66	164	222355	20.00	ng	-0.07
63) Phenanthrene-d10	11.15	188	358209	20.00	ng	-0.06
75) Chrysene-d12	13.80	240	255495	20.00	ng	-0.03
86) Perylene-d12	15.16	264	211020	20.00	ng	-0.05
System Monitoring Compounds						
5) 2-Fluorophenol	5.22	112	1051966	104.55	ng	-0.06
7) Phenol-d6	6.29	99	1345459	107.86	ng	-0.05
23) Nitrobenzene-d5	7.21	82	843453	83.58	ng	-0.06
41) 2,4,6-Tribromophenol	10.46	330	236675	102.04	ng	-0.06
44) 2-Fluorobiphenyl	9.00	172	1113473	83.98	ng	-0.06
78) Terphenyl-d14	12.75	244	643296	57.06	ng	-0.05
Target Compounds						Qvalue
9) Benzaldehyde	6.18	77	38268	5.20	ng	95
10) Phenol	6.30	94	75010	5.37	ng	89
17) 2-Methylphenol	6.91	107	40606	4.51	ng	98
20) 3+4-Methylphenols	7.07	107	319479	28.58	ng	90
27) 2,4-Dimethylphenol	7.60	122	109524	11.81	ng	92
49) Dimethylphthalate	9.40	163	252032	14.88	ng	# 91
70) Phenanthrene	11.17	178	86194	4.34	ng	# 88
74) Fluoranthene	12.36	202	75928	3.78	ng	72
77) Pyrene	12.59	202	86056	4.40	ng	# 73
91) Benzo(q,h,i)perylene	16.80	276	23818	2.33	ng	# 83

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

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