

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121416\
 Data File : BF091582.D
 Acq On : 14 Dec 2016 13:29
 Operator : UM/SJ
 Sample : H5961-01DL2 20X
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 QNWP8-MW26S-GWDL2

Quant Time: Dec 15 00:36:44 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 09 19:00:21 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	266594	20.00	ng	-0.02
21) Naphthalene-d8	8.09	136	1119476	20.00	ng	-0.01
38) Acenaphthene-d10	9.84	164	520191	20.00	ng	-0.02
63) Phenanthrene-d10	11.32	188	818262	20.00	ng	-0.01
75) Chrysene-d12	13.95	240	569758	20.00	ng	-0.01
86) Perylene-d12	15.37	264	530845	20.00	ng	-0.01
System Monitoring Compounds						
5) 2-Fluorophenol	5.39	112	42255	2.67	ng	-0.02
7) Phenol-d6	6.43	99	38395	1.95	ng	-0.02
23) Nitrobenzene-d5	7.36	82	79554	3.97	ng	-0.02
41) 2,4,6-Tribromophenol	10.62	330	30732	7.11	ng	-0.01
44) 2-Fluorobiphenyl	9.16	172	163592	5.43	ng	-0.02
78) Terphenyl-d14	12.91	244	83930	3.34	ng	-0.01
Target Compounds						Qvalue
27) 2,4-Dimethylphenol	7.74	122	63571	3.87	ng	94
31) Naphthalene	8.11	128	2624434	50.39	ng	99
37) 2-Methylnaphthalene	8.80	142	80586	2.39	ng	97
45) 1,1'-Biphenyl	9.26	154	127154	3.37	ng	94
51) Acenaphthene	9.87	154	289851	9.38	ng	96
54) Dibenzofuran	10.04	168	272622	7.13	ng	98
70) Phenanthrene	11.34	178	300375	7.41	ng	99
72) Carbazole	11.55	167	131327	3.43	ng	98

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

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