

Data Path : Z:\HPCHEM1\BNA_M\Data\BM052015\
 Data File : BM001375.D
 Acq On : 22 May 2015 00:42
 Operator : TP/IZ
 Sample : G2255-10
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 KY032MW006-150512

Quant Time: May 22 06:58:15 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM052015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 25 10:49:04 2015
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	0.000	152	0	0.00	ng	-7.43
6) Naphthalene-d8	10.148	136	7039	5.00	ng	#-0.05
12) Acenaphthene-d10	14.199	164	25551	5.00	ng	0.09
18) Phenanthrene-d10	16.859	188	79423	5.00	ng	# 0.00
25) Chrysene-d12	21.064	240	39221	5.00	ng	# 0.00
32) Perylene-d12	23.165	264	33209	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.067	112	9792	0.00	ng	0.05
5) Phenol-d6	6.567	99	1353	0.00	ng	-0.05
7) Nitrobenzene-d5	8.669	82	12844	33.83	ng	0.08
13) 2,4,6-Tribromophenol	15.636	330	10139	9.55	ng	0.04
14) 2-Fluorobiphenyl	12.872	172	73966	9.71	ng	0.15
27) Terphenyl-d14	19.513	244	50758	9.84	ng	0.00
Target Compounds						
8) Nitrobenzene	8.357	77	233802	485.35	ng	# 49
9) Naphthalene	10.319	128	44568	29.67	ng	# 4
10) Hexachlorobutadiene	10.576	225	83	0.29	ng	# 19
11) 2-Methylnaphthalene	11.893	142	834	0.78	ng	# 1
15) Acenaphthylene	13.945	152	5234	0.49	ng	# 33
16) Acenaphthene	14.272	154	516	0.08	ng	# 76
17) Fluorene	15.296	166	752	0.14	ng	# 14
19) 4-Bromophenyl-phenylether	15.636	248	1684	0.38	ng	# 1
20) Hexachlorobenzene	16.179	284	183	0.03	ng	# 1
21) Pentachlorophenol	16.519	266	88	0.42	ng	# 57
22) Phenanthrene	16.893	178	1777	0.06	ng	# 60
23) Anthracene	17.096	178	503	0.05	ng	# 26
24) Fluoranthene	18.919	202	5022	0.21	ng	# 76
26) Pyrene	19.283	202	462	0.04	ng	# 58
28) Benzo(a)anthracene	21.049	228	710	0.06	ng	# 13
29) Chrysene	21.049	228	691	0.06	ng	# 6
30) Bis(2-ethylhexyl)phtha...	20.987	149	160072	22.26	ng	99
31) Indeno(1,2,3-cd)pyrene	25.239	276	452	0.04	ng	# 66
33) Benzo(b)fluoranthene	22.539	252	479	0.05	ng	# 1
34) Benzo(k)fluoranthene	22.581	252	244	0.03	ng	# 1
35) Benzo(a)pyrene	23.071	252	324	0.04	ng	# 1
36) Dibenzo(a,h)anthracene	25.260	278	233	0.03	ng	# 1
37) Benzo(g,h,i)perylene	25.886	276	286	0.03	ng	# 1

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

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