

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071216\
 Data File : BG023036.D
 Acq On : 12 Jul 2016 14:44
 Operator : UM/SJ
 Sample : PB92009BL
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB92009BL

Quant Time: Jul 13 01:05:09 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 12 18:30:58 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	1335	20.00	ng	0.01
21) Naphthalene-d8	10.98	136	606	20.00	ng	0.00
38) Acenaphthene-d10	14.78	164	97	20.00	ng	-0.01
63) Phenanthrene-d10	17.66	188	8058	20.00	ng	0.10
75) Chrysene-d12	22.00	240	55	20.00	ng	0.15
86) Perylene-d12	25.38	264	358	20.00	ng	0.20
System Monitoring Compounds						
5) 2-Fluorophenol	5.70	112	778988	9543.38	ng	0.00
7) Phenol-d6	7.31	99	1254776	9814.76	ng	0.00
23) Nitrobenzene-d5	9.33	82	886474	62637.29	ng	0.00
41) 2,4,6-Tribromophenol	16.29	330	709035	406714.49	ng	0.00
44) 2-Fluorobiphenyl	13.42	172	1985887	322478.19	ng	0.00
78) Terphenyl-d14	20.16	244	2907731	-20.00	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.58	88	120	3.78	ng	# 73
3) Pyridine	3.99	79	725	6.68	ng	# 49
4) n-Nitrosodimethylamine	3.90	42	390	8.37	ng	# 1
6) Aniline	7.47	93	5830	32.91	ng	# 69
8) 2-Chlorophenol	7.71	128	573	6.60	ng	# 1
9) Benzaldehyde	7.31	77	2883	33.75	ng	# 18
10) Phenol	7.33	94	857	6.52	ng	# 1
11) bis(2-Chloroethyl)ether	7.58	93	327	3.14	ng	# 23
12) 1,3-Dichlorobenzene	8.04	146	543	5.11	ng	# 71
15) Benzyl Alcohol	8.47	79	15279	130.48	ng	# 73
16) 2,2'-oxybis(1-Chloropropan	8.67	45	473	3.71	ng	# 34
17) 2-Methylphenol	8.60	107	288	3.36	ng	# 68
18) Hexachloroethane	9.24	117	249	5.54	ng	# 1
22) Acetophenone	8.97	105	977	53.72	ng	# 52
24) Nitrobenzene	9.37	77	801	53.26	ng	# 50
25) Isophorone	9.88	82	430	15.11	ng	# 69
26) 2-Nitrophenol	10.09	139	92	15.59	ng	# 15
27) 2,4-Dimethylphenol	10.14	122	67	6.57	ng	# 48
28) bis(2-Chloroethoxy)methane	10.38	93	123	7.75	ng	# 50
29) 2,4-Dichlorophenol	10.65	162	302	27.76	ng	# 25
30) 1,2,4-Trichlorobenzene	10.85	180	112	8.98	ng	# 14
31) Naphthalene	11.04	128	1173	38.02	ng	# 67
32) Benzoic acid	10.34	122	79	8.51	ng	# 1
33) 4-Chloroaniline	11.15	127	4267	282.85	ng	# 81
34) Hexachlorobutadiene	11.31	225	172	17.02	ng	# 18
35) Caprolactam	11.99	113	66	14.75	ng	# 10
36) 4-Chloro-3-methylphenol	12.24	107	124	8.33	ng	# 26
37) 2-Methylnaphthalene	12.85	142	382	15.67	ng	# 96
39) 1,2,4,5-Tetrachlorobenzene	12.99	216	774	185.16	ng	# 34
40) Hexachlorocyclopentadiene	12.97	237	193	80.20	ng	# 26
42) 2,4,6-Trichlorophenol	13.24	196	431	179.39	ng	# 28
43) 2,4,5-Trichlorophenol	13.29	196	93	37.69	ng	# 32
45) 1,1'-Biphenyl	13.64	154	1068	146.74	ng	# 74
46) 2-Chloronaphthalene	13.69	162	1156	195.81	ng	# 31

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
47) 2-Nitroaniline	13.87	65	88	34.93	nq	# 41
48) Acenaphthylene	14.52	152	1005	113.48	nq	# 62
49) Dimethylphthalate	14.25	163	1178	148.67	nq	# 77
50) 2,6-Dinitrotoluene	14.37	165	161	90.41	nq	# 9
51) Acenaphthene	14.87	154	993	171.58	nq	# 40
52) 3-Nitroaniline	14.72	138	1042	586.83	nq	# 22
53) 2,4-Dinitrophenol	14.88	184	87	71.20	nq	# 9
54) Dibenzofuran	15.20	168	1414	163.23	nq	# 70
55) 4-Nitrophenol	15.07	139	263	176.70	nq	# 1
56) 2,4-Dinitrotoluene	15.14	165	79	30.43	nq	# 1
57) Fluorene	15.84	166	1258	168.73	nq	# 40
58) 2,3,4,6-Tetrachlorophenol	15.43	232	245	99.25	nq	# 75
59) Diethylphthalate	15.60	149	1269	157.40	nq	# 62
60) 4-Chlorophenyl-phenylether	15.83	204	595	131.03	nq	# 70
61) 4-Nitroaniline	15.89	138	232	120.75	nq	# 8
62) Azobenzene	16.16	77	458	59.39	nq	# 85
64) 4,6-Dinitro-2-methylphenol	16.10	198	150	2.50	nq	# 33
65) n-Nitrosodiphenylamine	16.29	169	28046	120.81	nq	# 40
68) Atrazine	17.00	200	1493	14.50	nq	# 89
70) Phenanthrene	17.68	178	1643	4.16	nq	# 57
74) Fluoranthene	19.60	202	1502	3.10	nq	# 85
76) Benzidine	20.16	184	74289	47114.57	nq	# 97
77) Pyrene	20.17	202	15478	5007.89	nq	# 71
79) Butylbenzylphthalate	20.96	149	509	415.78	nq	# 1
80) Benzo(a)anthracene	21.92	228	2136	694.15	nq	# 65
81) 3,3'-Dichlorobenzidine	21.84	252	217	160.66	nq	# 40
82) Chrysene	22.05	228	129	44.69	nq	# 13
83) Bis(2-ethylhexyl)phthalate	21.85	149	824	484.73	nq	# 48
84) Di-n-octyl phthalate	23.12	149	1087	383.43	nq	# 89
85) Indeno(1,2,3-cd)pyrene	29.22	276	222	60.32	nq	# 100
87) Benzo(b)fluoranthene	24.30	252	169	8.18	nq	# 1
88) Benzo(k)fluoranthene	24.37	252	231	11.79	nq	# 1
89) Benzo(a)pyrene	25.20	252	673	33.99	nq	# 31
90) Dibenzo(a,h)anthracene	29.31	278	306	15.57	nq	# 59
91) Benzo(a,h,i)perylene	30.47	276	349	17.73	nq	# 1

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

