

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG091915\
 Data File : BG018763.D
 Acq On : 18 Sep 2015 11:28
 Operator : TP
 Sample : PB85606BSD
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB85606BSD

Quant Time: Sep 18 23:55:07 2015
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG091115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 18 23:51:57 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.00	152	41258	20.00	ng	0.00
21) Naphthalene-d8	10.80	136	167970	20.00	ng	0.00
38) Acenaphthene-d10	14.63	164	116391	20.00	ng	0.00
63) Phenanthrene-d10	17.37	188	337421	20.00	ng	0.00
75) Chrysene-d12	21.63	240	354851	20.00	ng	0.00
86) Perylene-d12	24.82	264	360151	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.58	112	294300	123.22	ng	0.00
7) Phenol-d6	7.16	99	412110	120.34	ng	0.00
23) Nitrobenzene-d5	9.16	82	225856	75.24	ng	0.00
41) 2,4,6-Tribromophenol	16.12	330	218825	139.63	ng	0.00
44) 2-Fluorobiphenyl	13.25	172	632232	75.41	ng	0.00
78) Terphenyl-d14	19.98	244	1097375	81.21	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.47	88	29830	29.88	ng	# 100
3) Pyridine	3.86	79	90979	29.40	ng	97
4) n-Nitrosodimethylamine	3.78	42	32106	29.79	ng	# 67
6) Aniline	7.32	93	79137	16.77	ng	96
8) 2-Chlorophenol	7.57	128	108916	39.49	ng	94
9) Benzaldehyde	7.14	77	36135	19.60	ng	99
10) Phenol	7.19	94	146089	40.38	ng	96
11) bis(2-Chloroethyl)ether	7.42	93	101187	36.09	ng	98
12) 1,3-Dichlorobenzene	7.89	146	113253	35.30	ng	98
13) 1,4-Dichlorobenzene	8.03	146	116087	36.08	ng	98
14) 1,2-Dichlorobenzene	8.35	146	113407	36.91	ng	98
15) Benzyl Alcohol	8.23	79	92527	37.49	ng	97
16) 2,2'-oxybis(1-Chloropropan	8.53	45	110045	34.57	ng	97
17) 2-Methylphenol	8.45	107	97416	38.75	ng	96
18) Hexachloroethane	9.09	117	39371	34.22	ng	97
19) n-Nitroso-di-n-propylamine	8.80	70	76656	31.78	ng	96
20) 3+4-Methylphenols	8.77	107	141166	40.00	ng	96
22) Acetophenone	8.82	105	159469	37.48	ng	# 97
24) Nitrobenzene	9.20	77	115255	36.16	ng	99
25) Isophorone	9.72	82	222770	34.51	ng	100
26) 2-Nitrophenol	9.91	139	60220	40.50	ng	97
27) 2,4-Dimethylphenol	9.97	122	112945	41.82	ng	97
28) bis(2-Chloroethoxy)methane	10.20	93	144182	38.90	ng	97
29) 2,4-Dichlorophenol	10.45	162	121929	44.68	ng	96
30) 1,2,4-Trichlorobenzene	10.67	180	116097	38.39	ng	95
31) Naphthalene	10.85	128	341482	37.96	ng	98
32) Benzoic acid	10.10	122	79618	37.58	ng	96
33) 4-Chloroaniline	10.96	127	62956	15.47	ng	97
34) Hexachlorobutadiene	11.14	225	70514	39.55	ng	93
35) Caprolactam	11.72	113	47929	40.85	ng	91
36) 4-Chloro-3-methylphenol	12.08	107	133274	43.78	ng	94
37) 2-Methylnaphthalene	12.46	142	257968	39.18	ng	98
39) 1,2,4,5-Tetrachlorobenzene	12.82	216	136206	36.90	ng	99
40) Hexachlorocyclopentadiene	12.81	237	140020	75.53	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.06	196	107253	41.95	ng	97
43) 2,4,5-Trichlorophenol	13.14	196	123268	43.95	ng	97
45) 1,1'-Biphenyl	13.46	154	350422	37.86	ng	98
46) 2-Chloronaphthalene	13.50	162	273888	38.41	ng	98
47) 2-Nitroaniline	13.70	65	83651	38.98	ng	89
48) Acenaphthylene	14.35	152	487085	38.64	ng	98
49) Dimethylphthalate	14.08	163	374954	38.90	ng	98
50) 2,6-Dinitrotoluene	14.20	165	88716	42.36	ng	96
51) Acenaphthene	14.69	154	280301	38.22	ng	98
52) 3-Nitroaniline	14.53	138	60718	24.27	ng	86
53) 2,4-Dinitrophenol	14.74	184	56598	55.95	ng	96
54) Dibenzofuran	15.03	168	474456	41.63	ng	99
55) 4-Nitrophenol	14.84	139	161806	83.76	ng	93
56) 2,4-Dinitrotoluene	14.99	165	131693	44.34	ng	# 92
57) Fluorene	15.67	166	403594	41.13	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.26	232	115636	43.97	ng	# 97
59) Diethylphthalate	15.44	149	398469	39.64	ng	97
60) 4-Chlorophenyl-phenylether	15.67	204	193023	40.94	ng	95
61) 4-Nitroaniline	15.69	138	105612	39.15	ng	96
62) Azobenzene	15.95	77	358183	40.32	ng	98
64) 4,6-Dinitro-2-methylphenol	15.75	198	56574	33.01	ng	98
65) n-Nitrosodiphenylamine	15.88	169	359696	36.81	ng	98
66) 4-Bromophenyl-phenylether	16.56	248	134728	39.26	ng	96
67) Hexachlorobenzene	16.68	284	151302	40.58	ng	95
68) Atrazine	16.82	200	149876	38.57	ng	99
69) Pentachlorophenol	17.02	266	167535	72.86	ng	91
70) Phenanthrene	17.42	178	701917	39.55	ng	99
71) Anthracene	17.51	178	713864	39.72	ng	98
72) Carbazole	17.77	167	668341	38.42	ng	99
73) Di-n-butylphthalate	18.33	149	793008	38.65	ng	99
74) Fluoranthene	19.42	202	839752	38.20	ng	98
76) Benzidine	19.60	184	273844	25.70	ng	98
77) Pyrene	19.78	202	857575	39.32	ng	98
79) Butylbenzylphthalate	20.67	149	347855	39.24	ng	96
80) Benzo(a)anthracene	21.61	228	813979	39.84	ng	99
81) 3,3'-Dichlorobenzidine	21.52	252	199300	25.68	ng	96
82) Chrysene	21.68	228	732114	38.06	ng	99
83) Bis(2-ethylhexyl)phthalate	21.52	149	507058	39.65	ng	98
84) Di-n-octyl phthalate	22.71	149	866606	40.06	ng	99
85) Indeno(1,2,3-cd)pyrene	28.45	276	902832	36.96	ng	# 100
87) Benzo(b)fluoranthene	23.81	252	843256	40.38	ng	98
88) Benzo(k)fluoranthene	23.87	252	814835	38.94	ng	98
89) Benzo(a)pyrene	24.67	252	817546	41.14	ng	97
90) Dibenzo(a,h)anthracene	28.51	278	729309	37.27	ng	97
91) Benzo(g,h,i)perylene	29.58	276	735795	36.93	ng	97

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

