

Data Path : Z:\HPCHEM1\BNA G\DATA\BG120415\
 Data File : BG019939.D
 Acq On : 5 Dec 2015 13:09
 Operator : UM/NP
 Sample : PB86993BSD
 Misc :
 ALS Vial : 30 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB86993BSD

Manual Integrations
 APPROVED

Mmdadoda
 12/7/2015 6:29:49 PM

Quant Time: Dec 07 05:01:02 2015
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG120215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 02 18:50:01 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.12	152	28182	20.00	ng	-0.02
21) Naphthalene-d8	10.94	136	124970	20.00	ng	-0.02
38) Acenaphthene-d10	14.75	164	88984	20.00	ng	-0.02
63) Phenanthrene-d10	17.49	188	228147	20.00	ng	-0.02
75) Chrysene-d12	21.76	240	276231	20.00	ng	-0.03
86) Perylene-d12	25.04	264	258709	20.00	ng	-0.05

System Monitoring Compounds

5) 2-Fluorophenol	5.66	112	210611	135.07	ng	0.00
7) Phenol-d6	7.26	99	310231	131.77	ng	-0.01
23) Nitrobenzene-d5	9.28	82	201795	82.41	ng	-0.02
41) 2,4,6-Tribromophenol	16.23	330	160708	118.04	ng	-0.02
44) 2-Fluorobiphenyl	13.37	172	491251	75.38	ng	-0.02
78) Terphenyl-d14	20.09	244	885147	78.16	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.52	88	19151	31.53	ng	# 100
3) Pyridine	3.93	79	63537	34.44	ng	# 89
4) n-Nitrosodimethylamine	3.84	42	32885	33.89	ng	# 81
6) Aniline	7.43	93	78405	25.72	ng	# 81
8) 2-Chlorophenol	7.67	128	77248	40.63	ng	# 96
9) Benzaldehyde	7.25	77	28716	18.29	ng	# 92
10) Phenol	7.29	94	107365	43.33	ng	# 81
11) bis(2-Chloroethyl)ether	7.53	93	70163	38.23	ng	# 91
12) 1,3-Dichlorobenzene	8.00	146	79072	36.69	ng	# 90
13) 1,4-Dichlorobenzene	8.15	146	81118	36.44	ng	# 92
14) 1,2-Dichlorobenzene	8.47	146	77691	36.59	ng	# 92
15) Benzyl Alcohol	8.35	79	82450	39.91	ng	# 89
16) 2,2'-oxybis(1-Chloropropan	8.65	45	109581	36.61	ng	# 94
17) 2-Methylphenol	8.55	107	74180	42.41	ng	# 95
18) Hexachloroethane	9.21	117	30047	36.05	ng	# 92
19) n-Nitroso-di-n-propylamine	8.92	70	67024	36.10	ng	# 90
20) 3+4-Methylphenols	8.88	107	101509	41.21	ng	# 95
22) Acetophenone	8.94	105	118398	34.56	ng	# 92
24) Nitrobenzene	9.33	77	101331	38.15	ng	# 90
25) Isophorone	9.85	82	182568	34.78	ng	# 95
26) 2-Nitrophenol	10.04	139	42969	41.89	ng	# 78
27) 2,4-Dimethylphenol	10.09	122	85306	39.93	ng	# 92
28) bis(2-Chloroethoxy)methane	10.33	93	103461	40.33	ng	# 99
29) 2,4-Dichlorophenol	10.57	162	90958	41.68	ng	# 96
30) 1,2,4-Trichlorobenzene	10.79	180	88618	35.93	ng	# 97
31) Naphthalene	10.98	128	249418	37.26	ng	# 97
32) Benzoic acid	10.21	122	62771	44.15	ng	# 80
33) 4-Chloroaniline	11.08	127	51453	17.44	ng	# 92
34) Hexachlorobutadiene	11.27	225	61069	33.57	ng	# 99
35) Caprolactam	11.86	113	32615m	40.15	ng	# 95
36) 4-Chloro-3-methylphenol	12.19	107	107168	39.43	ng	# 96
37) 2-Methylnaphthalene	12.58	142	191932	39.12	ng	# 98
39) 1,2,4,5-Tetrachlorobenzene	12.94	216	118382	35.06	ng	# 99
40) Hexachlorocyclopentadiene	12.93	237	131471	68.28	ng	# 99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.18	196	83516	37.36	nq	97
43) 2,4,5-Trichlorophenol	13.25	196	93685	39.65	nq	95
45) 1,1'-Biphenyl	13.59	154	255312	34.96	nq	99
46) 2-Chloronaphthalene	13.63	162	204690	36.36	nq	95
47) 2-Nitroaniline	13.82	65	75406	42.41	nq	88
48) Acenaphthylene	14.47	152	344079	35.89	nq	100
49) Dimethylphthalate	14.20	163	286400	35.78	nq	98
50) 2,6-Dinitrotoluene	14.31	165	63655	42.30	nq	# 78
51) Acenaphthene	14.81	154	215433	37.03	nq	99
52) 3-Nitroaniline	14.64	138	47042	30.09	nq	96
53) 2,4-Dinitrophenol	14.84	184	65281	81.34	nq	# 83
54) Dibenzofuran	15.14	168	350825	40.53	nq	94
55) 4-Nitrophenol	14.94	139	112146	82.36	nq	# 75
56) 2,4-Dinitrotoluene	15.10	165	90720	43.21	nq	# 86
57) Fluorene	15.79	166	282181	36.43	nq	100
58) 2,3,4,6-Tetrachlorophenol	15.36	232	90662	41.46	nq	99
59) Diethylphthalate	15.55	149	304420	35.14	nq	97
60) 4-Chlorophenyl-phenylether	15.78	204	153087	34.42	nq	96
61) 4-Nitroaniline	15.81	138	68895	39.18	nq	# 77
62) Azobenzene	16.08	77	292755	40.61	nq	95
64) 4,6-Dinitro-2-methylphenol	15.86	198	53197	42.33	nq	94
65) n-Nitrosodiphenylamine	15.99	169	260547	39.46	nq	98
66) 4-Bromophenyl-phenylether	16.68	248	103021	36.34	nq	93
67) Hexachlorobenzene	16.79	284	111804	37.86	nq	95
68) Atrazine	16.93	200	111193	38.76	nq	98
69) Pentachlorophenol	17.13	266	144171	83.64	nq	96
70) Phenanthrene	17.53	178	496568	38.47	nq	97
71) Anthracene	17.62	178	502628	38.23	nq	96
72) Carbazole	17.89	167	447739	38.66	nq	97
73) Di-n-butylphthalate	18.44	149	539775	38.02	nq	98
74) Fluoranthene	19.53	202	621202	37.62	nq	95
76) Benzidine	19.71	184	183155	20.19	nq	97
77) Pyrene	19.89	202	641422	39.47	nq	97
79) Butylbenzylphthalate	20.77	149	250162	39.27	nq	99
80) Benzo(a)anthracene	21.74	228	647883	38.32	nq	99
81) 3,3'-Dichlorobenzidine	21.65	252	146063	22.68	nq	98
82) Chrysene	21.81	228	592294	36.89	nq	96
83) Bis(2-ethylhexyl)phthalate	21.65	149	363703	38.91	nq	97
84) Di-n-octyl phthalate	22.89	149	621609	40.90	nq	# 94
85) Indeno(1,2,3-cd)pyrene	28.80	276	725658	41.03	nq	# 100
87) Benzo(b)fluoranthene	24.00	252	626194	38.19	nq	# 96
88) Benzo(k)fluoranthene	24.07	252	619878	38.17	nq	# 95
89) Benzo(a)pyrene	24.88	252	602943	38.73	nq	# 96
90) Dibenzo(a,h)anthracene	28.87	278	594626	38.66	nq	# 94
91) Benzo(g,h,i)perylene	29.97	276	583016	38.61	nq	# 93

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

