

Data Path : Z:\HPCHEM1\BNA G\DATA\BG060717\
 Data File : BG027290.D
 Acq On : 7 Jun 2017 19:18
 Operator : SJ/MA
 Sample : PB99548BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB99548BS

Manual Integrations
 APPROVED

mohammad
 6/8/2017 2:22:58 PM

Quant Time: Jun 08 06:50:26 2017
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 05 16:00:35 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.56	152	54578	20.00	ng	-0.02
21) Naphthalene-d8	11.38	136	274222	20.00	ng	-0.02
38) Acenaphthene-d10	15.14	164	185403	20.00	ng	-0.02
63) Phenanthrene-d10	17.89	188	416830	20.00	ng	-0.03
75) Chrysene-d12	22.28	240	443204	20.00	ng	-0.03
86) Perylene-d12	25.97	264	407561	20.00	ng	-0.06

System Monitoring Compounds

5) 2-Fluorophenol	6.14	112	518993	145.11	ng	-0.02
7) Phenol-d6	7.69	99	756561	144.10	ng	0.00
23) Nitrobenzene-d5	9.73	82	414315	95.02	ng	-0.02
41) 2,4,6-Tribromophenol	16.62	330	338465	138.63	ng	-0.02
44) 2-Fluorobiphenyl	13.77	172	1062038	80.14	ng	-0.02
78) Terphenyl-d14	20.47	244	1452729	83.62	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	4.16	88	42933	28.95	ng	# 83
3) Pyridine	4.53	79	135657	29.67	ng	# 73
4) n-Nitrosodimethylamine	4.44	42	67195	39.70	ng	# 46
6) Aniline	7.89	93	241256	36.49	ng	# 93
8) 2-Chlorophenol	8.13	128	173186	45.86	ng	92
9) Benzaldehyde	7.71	77	82809	22.81	ng	87
10) Phenol	7.72	94	259956	48.42	ng	# 76
11) bis(2-Chloroethyl)ether	7.96	93	169242	38.42	ng	87
12) 1,3-Dichlorobenzene	8.45	146	154941m	36.37	ng	
13) 1,4-Dichlorobenzene	8.59	146	161506	36.94	ng	96
14) 1,2-Dichlorobenzene	8.92	146	155260	36.41	ng	93
15) Benzyl Alcohol	8.79	79	175858	47.54	ng	# 83
16) 2,2'-oxybis(1-Chloropropan	9.07	45	266213	42.24	ng	93
17) 2-Methylphenol	8.97	107	173563	45.73	ng	# 87
18) Hexachloroethane	9.65	117	54657	37.12	ng	87
19) n-Nitroso-di-n-propylamine	9.35	70	149735	40.77	ng	# 85
20) 3+4-Methylphenols	9.30	107	240620	45.77	ng	88
22) Acetophenone	9.37	105	275191	39.24	ng	# 84
24) Nitrobenzene	9.77	77	210026	42.28	ng	# 86
25) Isophorone	10.29	82	427385	41.68	ng	# 96
26) 2-Nitrophenol	10.48	139	89538	44.36	ng	# 77
27) 2,4-Dimethylphenol	10.51	122	205079	46.52	ng	93
28) bis(2-Chloroethoxy)methane	10.76	93	266772	40.18	ng	100
29) 2,4-Dichlorophenol	11.01	162	185153	46.01	ng	96
30) 1,2,4-Trichlorobenzene	11.24	180	165764	37.20	ng	98
31) Naphthalene	11.43	128	556691	37.14	ng	99
32) Benzoic acid	10.63	122	118039	40.26	ng	# 86
33) 4-Chloroaniline	11.52	127	140081	22.43	ng	# 89
34) Hexachlorobutadiene	11.70	225	96355	35.18	ng	98
35) Caprolactam	12.30	113	71527m	51.82	ng	
36) 4-Chloro-3-methylphenol	12.59	107	236456	46.88	ng	95
37) 2-Methylnaphthalene	12.99	142	428421m	39.19	ng	
39) 1,2,4,5-Tetrachlorobenzene	13.35	216	208646	36.12	ng	98
40) Hexachlorocyclopentadiene	13.33	237	274146	89.18	ng	96

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42) 2,4,6-Trichlorophenol	13.58	196	153506	43.82	nq	99
43) 2,4,5-Trichlorophenol	13.65	196	171035	43.63	nq	95
45) 1,1'-Biphenyl	13.97	154	576293	38.08	nq	98
46) 2-Chloronaphthalene	14.03	162	437349	38.55	nq	99
47) 2-Nitroaniline	14.22	65	153112	47.53	nq	# 87
48) Acenaphthylene	14.87	152	714156	37.19	nq	98
49) Dimethylphthalate	14.58	163	605871	41.19	nq	98
50) 2,6-Dinitrotoluene	14.70	165	126587	43.03	nq	# 78
51) Acenaphthene	15.21	154	500588	40.08	nq	99
52) 3-Nitroaniline	15.03	138	97475	33.11	nq	91
53) 2,4-Dinitrophenol	15.23	184	101406	73.28	nq	96
54) Dibenzofuran	15.54	168	716617	41.06	nq	92
55) 4-Nitrophenol	15.31	139	214122	77.30	nq	# 58
56) 2,4-Dinitrotoluene	15.48	165	177602	46.69	nq	# 86
57) Fluorene	16.18	166	585495	37.77	nq	100
58) 2,3,4,6-Tetrachlorophenol	15.75	232	160273	45.80	nq	98
59) Diethylphthalate	15.93	149	606815	41.36	nq	97
60) 4-Chlorophenyl-phenylether	16.17	204	302087	38.68	nq	96
61) 4-Nitroaniline	16.20	138	133547	42.83	nq	# 67
62) Azobenzene	16.46	77	618626	45.50	nq	90
64) 4,6-Dinitro-2-methylphenol	16.24	198	85193	46.55	nq	87
65) n-Nitrosodiphenylamine	16.38	169	524648	40.15	nq	99
66) 4-Bromophenyl-phenylether	17.07	248	200334	40.61	nq	97
67) Hexachlorobenzene	17.19	284	212296	40.01	nq	91
68) Atrazine	17.32	200	185230	42.39	nq	96
69) Pentachlorophenol	17.53	266	240017	77.17	nq	97
70) Phenanthrene	17.93	178	929300	39.63	nq	96
71) Anthracene	18.02	178	928504	39.52	nq	98
72) Carbazole	18.29	167	791414	35.71	nq	95
73) Di-n-butylphthalate	18.82	149	922733	42.94	nq	# 93
74) Fluoranthene	19.92	202	973690	38.86	nq	94
76) Benzidine	20.09	184	344849	25.83	nq	97
77) Pyrene	20.29	202	1000988	37.38	nq	95
79) Butylbenzylphthalate	21.17	149	397446	42.00	nq	93
80) Benzo(a)anthracene	22.25	228	1000925	39.60	nq	96
81) 3,3'-Dichlorobenzidine	22.15	252	257744	26.21	nq	# 98
82) Chrysene	22.33	228	930944	38.36	nq	96
83) Bis(2-ethylhexyl)phthalate	22.12	149	574148	40.15	nq	99
84) Di-n-octyl phthalate	23.50	149	980706	41.41	nq	# 89
85) Indeno(1,2,3-cd)pyrene	30.22	276	1157802	39.40	nq	# 89
87) Benzo(b)fluoranthene	24.79	252	988487	39.11	nq	# 96
88) Benzo(k)fluoranthene	24.87	252	976787	39.39	nq	# 93
89) Benzo(a)pyrene	25.79	252	960718	40.38	nq	# 94
90) Dibenzo(a,h)anthracene	30.30	278	999059	40.11	nq	# 94
91) Benzo(g,h,i)perylene	31.56	276	943292	39.10	nq	# 90

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

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