

Data Path : U:\HPCHEM1\BNA G\DATA\BG020218\
 Data File : BG032519.D
 Acq On : 2 Feb 2018 19:31
 Operator : SJ/JU
 Sample : PB106268BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB106268BS

Manual Integrations
 APPROVED

Sohil
 2/5/2018 11:15:13 AM

Quant Time: Feb 05 10:08:05 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG012918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 29 15:32:55 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.71	152	24125	20.00	ng	-0.01
21) Naphthalene-d8	11.59	136	95161	20.00	ng	-0.01
38) Acenaphthene-d10	15.34	164	72451	20.00	ng	-0.01
63) Phenanthrene-d10	18.08	188	200029	20.00	ng	-0.01
75) Chrysene-d12	22.41	240	260169	20.00	ng	-0.01
86) Perylene-d12	26.08	264	283984	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	6.15	112	148873	123.95	ng	0.00
7) Phenol-d6	7.83	99	194905	121.60	ng	0.00
23) Nitrobenzene-d5	9.91	82	121842	79.97	ng	-0.01
41) 2,4,6-Tribromophenol	16.82	330	180079	130.58	ng	-0.01
44) 2-Fluorobiphenyl	13.97	172	471258	77.31	ng	-0.02
78) Terphenyl-d14	20.61	244	1047372	86.23	ng	-0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.80	88	13484	33.366	ng	# 73
3) Pyridine	4.25	79	34846	31.934	ng	# 85
4) n-Nitrosodimethylamine	4.17	42	23644	41.624	ng	# 64
6) Aniline	8.01	93	24478	12.249	ng	94
8) 2-Chlorophenol	8.26	128	60515	42.623	ng	85
9) Benzaldehyde	7.81	77	17699	21.380	ng	92
10) Phenol	7.86	94	64958	42.677	ng	84
11) bis(2-Chloroethyl)ether	8.10	93	44698	38.541	ng	# 81
12) 1,3-Dichlorobenzene	8.60	146	71863	37.427	ng	90
13) 1,4-Dichlorobenzene	8.75	146	73597	38.243	ng	91
14) 1,2-Dichlorobenzene	9.08	146	71234	37.785	ng	98
15) Benzyl Alcohol	8.95	79	49728	37.646	ng	98
16) 2,2'-oxybis(1-Chloropropan	9.23	45	38979	35.469	ng	71
17) 2-Methylphenol	9.16	107	49078	39.701	ng	# 94
18) Hexachloroethane	9.82	117	23463	36.313	ng	87
19) n-Nitroso-di-n-propylamine	9.52	70	36274	34.133	ng	# 90
20) 3+4-Methylphenols	9.49	107	63943	36.875	ng	92
22) Acetophenone	9.55	105	82709	37.078	ng	# 89
24) Nitrobenzene	9.95	77	60403	41.000	ng	# 87
25) Isophorone	10.47	82	108847	41.958	ng	# 84
26) 2-Nitrophenol	10.68	139	36818	41.342	ng	# 76
27) 2,4-Dimethylphenol	10.72	122	59667	46.679	ng	99
28) bis(2-Chloroethoxy)methane	10.96	93	59216	41.627	ng	# 95
29) 2,4-Dichlorophenol	11.22	162	78785	46.527	ng	83
30) 1,2,4-Trichlorobenzene	11.44	180	86316	38.246	ng	96
31) Naphthalene	11.64	128	185919	40.002	ng	97
32) Benzoic acid	10.81	122	26114m	30.514	ng	
33) 4-Chloroaniline	11.74	127	27025	13.037	ng	# 89
34) Hexachlorobutadiene	11.91	225	66403	37.817	ng	96
35) Caprolactam	12.49	113	22305	45.885	ng	# 48
36) 4-Chloro-3-methylphenol	12.82	107	72367	45.055	ng	# 86
37) 2-Methylnaphthalene	13.21	142	160377	41.074	ng	92
39) 1,2,4,5-Tetrachlorobenzene	13.56	216	127617	37.651	ng	98
40) Hexachlorocyclopentadiene	13.53	237	167341	79.018	ng	92

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.79	196	78529	42.594	nq	98
43) 2,4,5-Trichlorophenol	13.87	196	92048	42.829	nq	# 93
45) 1,1'-Biphenyl	14.18	154	235788	38.521	nq	95
46) 2-Chloronaphthalene	14.24	162	187511	39.085	nq	96
47) 2-Nitroaniline	14.43	65	45077	42.488	nq	# 87
48) Acenaphthylene	15.07	152	283718	39.119	nq	97
49) Dimethylphthalate	14.77	163	263328	39.419	nq	98
50) 2,6-Dinitrotoluene	14.91	165	57359	40.909	nq	# 75
51) Acenaphthene	15.41	154	185314	36.842	nq	99
52) 3-Nitroaniline	15.24	138	24776	20.230	nq	# 74
53) 2,4-Dinitrophenol	15.44	184	11572	16.648	nq	# 91
54) Dibenzofuran	15.74	168	307908	41.359	nq	97
55) 4-Nitrophenol	15.54	139	70694	77.104	nq	# 74
56) 2,4-Dinitrotoluene	15.69	165	87669	43.915	nq	# 73
57) Fluorene	16.38	166	253271	40.940	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.96	232	95620	45.109	nq	# 87
59) Diethylphthalate	16.12	149	266685	40.989	nq	95
60) 4-Chlorophenyl-phenylether	16.36	204	163115	41.813	nq	95
61) 4-Nitroaniline	16.40	138	51366	39.132	nq	# 72
62) Azobenzene	16.65	77	159357	40.985	nq	81
64) 4,6-Dinitro-2-methylphenol	16.45	198	23362m	15.634	nq	
65) n-Nitrosodiphenylamine	16.57	169	238645	39.977	nq	99
66) 4-Bromophenyl-phenylether	17.25	248	120445	40.644	nq	94
67) Hexachlorobenzene	17.38	284	129276	39.420	nq	# 92
68) Atrazine	17.50	200	114146	44.543	nq	96
69) Pentachlorophenol	17.72	266	120248	58.532	nq	96
70) Phenanthrene	18.12	178	456406	40.915	nq	99
71) Anthracene	18.21	178	469070	42.231	nq	99
72) Carbazole	18.48	167	403338	41.099	nq	98
73) Di-n-butylphthalate	18.98	149	462836	42.597	nq	99
74) Fluoranthene	20.09	202	624725	42.705	nq	95
76) Benzidine	20.25	184	151140	29.585	nq	99
77) Pyrene	20.44	202	627582	39.329	nq	99
79) Butylbenzylphthalate	21.30	149	216621	43.038	nq	# 74
80) Benzo(a)anthracene	22.39	228	677946	42.033	nq	99
81) 3,3'-Dichlorobenzidine	22.28	252	113584	18.176	nq	# 97
82) Chrysene	22.46	228	630367	40.469	nq	96
83) Bis(2-ethylhexyl)phthalate	22.22	149	302871	42.438	nq	# 97
84) Di-n-octyl phthalate	23.59	149	528796	46.715	nq	# 89
85) Indeno(1,2,3-cd)pyrene	30.34	276	860224	43.656	nq	# 84
87) Benzo(b)fluoranthene	24.90	252	738991	43.067	nq	# 95
88) Benzo(k)fluoranthene	24.98	252	701550	41.272	nq	# 96
89) Benzo(a)pyrene	25.91	252	711292	43.481	nq	# 94
90) Dibenzo(a,h)anthracene	30.41	278	707139	42.257	nq	# 95
91) Benzo(g,h,i)perylene	31.68	276	711873	42.857	nq	# 90

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

