

Data Path : Z:\HPCHEM1\BNA G\DATA\BG021916\
 Data File : BG020985.D
 Acq On : 19 Feb 2016 23:32
 Operator : SJ/UM
 Sample : PB88453BSD
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PB88453BSD

Manual Integrations
 APPROVED

UMANGI
 2/22/2016 10:55:38 AM

Quant Time: Feb 20 02:32:30 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 11 15:26:53 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.26	152	19238	20.00	ng	0.00
21) Naphthalene-d8	11.09	136	98717	20.00	ng	0.00
38) Acenaphthene-d10	14.87	164	63356	20.00	ng	-0.02
63) Phenanthrene-d10	17.61	188	138466	20.00	ng	0.00
75) Chrysene-d12	21.89	240	129167	20.00	ng	-0.02
86) Perylene-d12	25.26	264	122202	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.80	112	107435	94.05	ng	0.00
7) Phenol-d6	7.41	99	156751	93.88	ng	0.00
23) Nitrobenzene-d5	9.43	82	125127	83.34	ng	0.00
41) 2,4,6-Tribromophenol	16.35	330	62249	99.91	ng	0.00
44) 2-Fluorobiphenyl	13.50	172	360337	79.89	ng	0.00
78) Terphenyl-d14	20.19	244	491162	88.71	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.63	88	13281	31.97	ng	93
3) Pyridine	4.04	79	41117	33.47	ng	98
4) n-Nitrosodimethylamine	3.95	42	12084	37.59	ng	94
6) Aniline	7.58	93	32438	15.58	ng	100
8) 2-Chlorophenol	7.82	128	64522	45.21	ng	99
9) Benzaldehyde	7.39	77	8670	10.37	ng	97
10) Phenol	7.43	94	81162	45.93	ng	98
11) bis(2-Chloroethyl)ether	7.67	93	55280	39.17	ng	94
12) 1,3-Dichlorobenzene	8.15	146	59230	39.26	ng	100
13) 1,4-Dichlorobenzene	8.30	146	61576	40.25	ng	98
14) 1,2-Dichlorobenzene	8.62	146	60162	40.43	ng	99
15) Benzyl Alcohol	8.50	79	46929	44.59	ng	96
16) 2,2'-oxybis(1-Chloropropan	8.79	45	53707	36.48	ng	98
17) 2-Methylphenol	8.70	107	57446	45.13	ng	97
18) Hexachloroethane	9.36	117	21313	40.52	ng	97
19) n-Nitroso-di-n-propylamine	9.06	70	41820	38.82	ng	# 96
20) 3+4-Methylphenols	9.03	107	80751	45.52	ng	94
22) Acetophenone	9.09	105	88942	37.38	ng	# 99
24) Nitrobenzene	9.47	77	61017	38.99	ng	97
25) Isophorone	10.00	82	123428	39.00	ng	98
26) 2-Nitrophenol	10.19	139	35977	43.68	ng	97
27) 2,4-Dimethylphenol	10.24	122	68636	43.49	ng	99
28) bis(2-Chloroethoxy)methane	10.47	93	81022	41.55	ng	99
29) 2,4-Dichlorophenol	10.72	162	68383	48.34	ng	98
30) 1,2,4-Trichlorobenzene	10.94	180	58146	40.81	ng	98
31) Naphthalene	11.14	128	208552	40.84	ng	100
32) Benzoic acid	10.37	122	51267	40.50	ng	100
33) 4-Chloroaniline	11.24	127	15560	7.08	ng	98
34) Hexachlorobutadiene	11.41	225	29190	40.19	ng	98
35) Caprolactam	12.01	113	27631m	50.98	ng	
36) 4-Chloro-3-methylphenol	12.34	107	79369	48.97	ng	99
37) 2-Methylnaphthalene	12.72	142	164086	45.79	ng	100
39) 1,2,4,5-Tetrachlorobenzene	13.08	216	68825	37.03	ng	98
40) Hexachlorocyclopentadiene	13.06	237	63387	77.52	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.31	196	56802	45.50	nq	99
43) 2,4,5-Trichlorophenol	13.39	196	63081	48.06	nq	99
45) 1,1'-Biphenyl	13.71	154	214389	38.36	nq	99
46) 2-Chloronaphthalene	13.76	162	162537	40.62	nq	99
47) 2-Nitroaniline	13.96	65	43485	45.38	nq	98
48) Acenaphthylene	14.60	152	295513	42.47	nq	99
49) Dimethylphthalate	14.32	163	219635	44.33	nq	99
50) 2,6-Dinitrotoluene	14.44	165	49335	46.47	nq	98
51) Acenaphthene	14.94	154	177767	42.14	nq	98
52) 3-Nitroaniline	14.77	138	24247	19.66	nq	97
53) 2,4-Dinitrophenol	14.98	184	47562	102.95	nq	89
54) Dibenzofuran	15.27	168	269739	48.34	nq	99
55) 4-Nitrophenol	15.07	139	93298	100.15	nq	98
56) 2,4-Dinitrotoluene	15.22	165	69443	50.60	nq	95
57) Fluorene	15.91	166	233822	45.50	nq	99
58) 2,3,4,6-Tetrachlorophenol	15.49	232	54331	54.32	nq	99
59) Diethylphthalate	15.67	149	236322	46.40	nq	99
60) 4-Chlorophenyl-phenylether	15.90	204	101831	44.62	nq	98
61) 4-Nitroaniline	15.93	138	56890	45.94	nq	99
62) Azobenzene	16.19	77	193119	48.23	nq	98
64) 4,6-Dinitro-2-methylphenol	15.98	198	31292	44.88	nq	98
65) n-Nitrosodiphenylamine	16.11	169	198628	44.40	nq	98
66) 4-Bromophenyl-phenylether	16.79	248	63165	43.19	nq	99
67) Hexachlorobenzene	16.92	284	68961	44.31	nq	97
68) Atrazine	17.05	200	60980	44.92	nq	99
69) Pentachlorophenol	17.26	266	81243	91.07	nq	97
70) Phenanthrene	17.65	178	347816	44.12	nq	99
71) Anthracene	17.74	178	344032	43.00	nq	100
72) Carbazole	18.01	167	321779	44.81	nq	100
73) Di-n-butylphthalate	18.54	149	394075	42.80	nq	99
74) Fluoranthene	19.64	202	365729	42.95	nq	98
76) Benzidine	19.81	184	96652	23.14	nq	97
77) Pyrene	20.00	202	378074	44.37	nq	99
79) Butylbenzylphthalate	20.87	149	179015	43.89	nq	99
80) Benzo(a)anthracene	21.87	228	345840	44.85	nq	100
81) 3,3'-Dichlorobenzidine	21.77	252	44879	15.68	nq	98
82) Chrysene	21.94	228	304090	41.93	nq	100
83) Bis(2-ethylhexyl)phthalate	21.75	149	261182	43.19	nq	100
84) Di-n-octyl phthalate	23.01	149	435753	43.90	nq	99
85) Indeno(1,2,3-cd)pyrene	29.12	276	368693	44.21	nq	98
87) Benzo(b)fluoranthene	24.18	252	335463	44.86	nq	99
88) Benzo(k)fluoranthene	24.25	252	315251	43.02	nq	99
89) Benzo(a)pyrene	25.10	252	316856	44.47	nq	100
90) Dibenzo(a,h)anthracene	29.19	278	301321	43.49	nq	99
91) Benzo(g,h,i)perylene	30.33	276	298847	43.45	nq	98

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

