

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122216\
 Data File : BG025391.D
 Acq On : 22 Dec 2016 16:22
 Operator : UM/SJ
 Sample : H6103-06
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 LOQ-WATER-06-QT1-2017

Quant Time: Dec 23 00:28:56 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 21 18:42:01 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.22	152	51603	20.00	ng	0.00
21) Naphthalene-d8	11.04	136	206127	20.00	ng	0.00
38) Acenaphthene-d10	14.84	164	164637	20.00	ng	0.00
63) Phenanthrene-d10	17.59	188	390376	20.00	ng	0.00
75) Chrysene-d12	21.88	240	571352	20.00	ng	0.00
86) Perylene-d12	25.29	264	589723	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.75	112	477624	168.24	ng	0.00
7) Phenol-d6	7.38	99	649990	162.57	ng	0.00
23) Nitrobenzene-d5	9.40	82	529213	113.42	ng	0.00
41) 2,4,6-Tribromophenol	16.33	330	408464	154.10	ng	0.00
44) 2-Fluorobiphenyl	13.46	172	1127459	110.87	ng	0.00
78) Terphenyl-d14	20.17	244	2212934	102.69	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	3.59	88	19400	16.02	ng	# 92
3) Pyridine	4.00	79	56517	16.10	ng	94
4) n-Nitrosodimethylamine	3.92	42	33733	16.18	ng	# 97
6) Aniline	7.53	93	57567	10.63	ng	96
8) 2-Chlorophenol	7.78	128	57125	17.84	ng	93
9) Benzaldehyde	7.36	77	52227	17.85	ng	86
10) Phenol	7.41	94	82027	18.51	ng	94
11) bis(2-Chloroethyl)ether	7.63	93	56436	16.52	ng	97
12) 1,3-Dichlorobenzene	8.10	146	64733	16.70	ng	95
13) 1,4-Dichlorobenzene	8.25	146	68102	16.96	ng	94
14) 1,2-Dichlorobenzene	8.57	146	62556	16.32	ng	95
15) Benzyl Alcohol	8.46	79	69049	18.09	ng	95
16) 2,2'-oxybis(1-Chloropropan	8.73	45	108011	17.08	ng	93
17) 2-Methylphenol	8.67	107	53865	18.51	ng	97
18) Hexachloroethane	9.30	117	25965	16.83	ng	85
19) n-Nitroso-di-n-propylamine	9.01	70	54645	16.18	ng	98
20) 3+4-Methylphenols	9.00	107	70229	17.43	ng	89
22) Acetophenone	9.04	105	98230	17.15	ng	# 93
24) Nitrobenzene	9.44	77	91434	17.87	ng	97
25) Isophorone	9.96	82	142253	16.84	ng	97
26) 2-Nitrophenol	10.16	139	35983	18.38	ng	97
27) 2,4-Dimethylphenol	10.20	122	61605	19.14	ng	94
28) bis(2-Chloroethoxy)methane	10.43	93	76362	17.41	ng	94
29) 2,4-Dichlorophenol	10.70	162	62767	18.23	ng	93
30) 1,2,4-Trichlorobenzene	10.90	180	73273	17.61	ng	96
31) Naphthalene	11.09	128	179182	17.41	ng	96
32) Benzoic acid	10.32	122	44683	17.27	ng	92
33) 4-Chloroaniline	11.22	127	38247	8.83	ng	# 85
34) Hexachlorobutadiene	11.35	225	57914	17.05	ng	96
35) Caprolactam	11.98	113	22211	17.17	ng	# 82
36) 4-Chloro-3-methylphenol	12.32	107	76725	18.05	ng	95
37) 2-Methylnaphthalene	12.68	142	137760	18.06	ng	91
39) 1,2,4,5-Tetrachlorobenzene	13.05	216	96383	17.73	ng	96
40) Hexachlorocyclopentadiene	13.00	237	17633	16.45	ng	89

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.29	196	61500	17.97	nq	94
43) 2,4,5-Trichlorophenol	13.37	196	65441	18.27	nq #	89
45) 1,1'-Biphenyl	13.67	154	195624	17.67	nq	96
46) 2-Chloronaphthalene	13.73	162	152023	17.57	nq	97
47) 2-Nitroaniline	13.94	65	65229	17.86	nq	95
48) Acenaphthylene	14.57	152	238286	17.77	nq	94
49) Dimethylphthalate	14.28	163	212165	17.68	nq #	95
50) 2,6-Dinitrotoluene	14.42	165	45374	17.30	nq	98
51) Acenaphthene	14.90	154	150775	17.19	nq	97
52) 3-Nitroaniline	14.76	138	34387	13.64	nq	89
53) 2,4-Dinitrophenol	15.00	184	11408	11.29	nq #	90
54) Dibenzofuran	15.24	168	247047	17.82	nq	98
55) 4-Nitrophenol	15.10	139	31629	16.80	nq #	78
56) 2,4-Dinitrotoluene	15.22	165	67933	17.79	nq #	91
57) Fluorene	15.88	166	206582	17.80	nq	95
58) 2,3,4,6-Tetrachlorophenol	15.47	232	71817	18.71	nq	93
59) Diethylphthalate	15.63	149	215297	17.11	nq	99
60) 4-Chlorophenyl-phenylether	15.87	204	115104	17.50	nq	91
61) 4-Nitroaniline	15.93	138	45133	17.79	nq	95
62) Azobenzene	16.16	77	216431	17.83	nq	97
64) 4,6-Dinitro-2-methylphenol	15.98	198	41766	15.45	nq	86
65) n-Nitrosodiphenylamine	16.08	169	184534	17.49	nq	95
66) 4-Bromophenyl-phenylether	16.76	248	81598	16.94	nq	98
67) Hexachlorobenzene	16.88	284	92586	16.75	nq	94
68) Atrazine	17.02	200	63063	13.64	nq	96
69) Pentachlorophenol	17.25	266	39789	13.89	nq	98
70) Phenanthrene	17.63	178	350472m	17.42	nq	
71) Anthracene	17.72	178	357951	17.51	nq	97
72) Carbazole	18.00	167	332379	18.18	nq #	94
73) Di-n-butylphthalate	18.50	149	405979	18.05	nq	99
74) Fluoranthene	19.63	202	494256	18.60	nq	90
76) Benzidine	19.80	184	135052	9.48	nq	97
77) Pyrene	19.99	202	521537	17.47	nq	93
79) Butylbenzylphthalate	20.84	149	207445	17.30	nq	94
80) Benzo(a)anthracene	21.86	228	571506	17.93	nq	91
81) 3,3'-Dichlorobenzidine	21.76	252	137602	11.12	nq	95
82) Chrysene	21.93	228	545756	17.86	nq	95
83) Bis(2-ethylhexyl)phthalate	21.70	149	290851	17.43	nq	98
84) Di-n-octyl phthalate	22.96	149	497056	17.94	nq #	93
85) Indeno(1,2,3-cd)pyrene	29.21	276	670829	17.27	nq #	96
87) Benzo(b)fluoranthene	24.19	252	585352	18.19	nq #	95
88) Benzo(k)fluoranthene	24.26	252	549697m	17.69	nq	
89) Benzo(a)pyrene	25.13	252	556931	17.92	nq #	97
90) Dibenzo(a,h)anthracene	29.26	278	584904	17.76	nq	98
91) Benzo(g,h,i)perylene	30.45	276	556127	16.99	nq	93

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

