

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062116\
 Data File : BG022480.D
 Acq On : 22 Jun 2016 00:14
 Operator : UM/SJ
 Sample : H3600-01MS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 PM-STA-1363(0-4)MS

Quant Time: Jun 22 13:51:13 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060616.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 22 02:15:22 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.98	152	21435	20.00	ng	0.00
21) Naphthalene-d8	10.80	136	111943	20.00	ng	0.00
38) Acenaphthene-d10	14.63	164	69184	20.00	ng	0.00
63) Phenanthrene-d10	17.38	188	153260	20.00	ng	0.00
75) Chrysene-d12	21.64	240	132750	20.00	ng	0.00
86) Perylene-d12	24.84	264	118521	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.54	112	186830	168.52	ng	0.00
7) Phenol-d6	7.17	99	265620	152.39	ng	0.00
23) Nitrobenzene-d5	9.16	82	163401	101.77	ng	0.00
41) 2,4,6-Tribromophenol	16.13	330	91020	116.43	ng	0.00
44) 2-Fluorobiphenyl	13.25	172	432255	95.21	ng	0.00
78) Terphenyl-d14	19.98	244	567218	101.44	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	3.33	88	18299	34.42	ng	# 64
3) Pyridine	3.76	79	42042	29.27	ng	99
4) n-Nitrosodimethylamine	3.67	42	15922	49.57	ng	94
6) Aniline	7.31	93	72530m	32.75	ng	
8) 2-Chlorophenol	7.55	128	77639	50.67	ng	# 79
10) Phenol	7.20	94	88942	49.49	ng	80
11) bis(2-Chloroethyl)ether	7.40	93	70481	49.19	ng	# 74
12) 1,3-Dichlorobenzene	7.87	146	65007	39.58	ng	# 91
13) 1,4-Dichlorobenzene	8.02	146	67697	41.37	ng	93
14) 1,2-Dichlorobenzene	8.34	146	68860	41.74	ng	93
15) Benzyl Alcohol	8.24	79	56082	49.80	ng	# 89
16) 2,2'-oxybis(1-Chloropropan	8.52	45	75908	51.06	ng	80
17) 2-Methylphenol	8.46	107	69341	51.71	ng	# 84
18) Hexachloroethane	9.06	117	25738	43.07	ng	84
19) n-Nitroso-di-n-propylamine	8.79	70	58822	49.85	ng	# 72
20) 3+4-Methylphenols	8.79	107	95038	49.41	ng	93
22) Acetophenone	8.82	105	116575	48.32	ng	# 83
24) Nitrobenzene	9.20	77	81729	50.62	ng	# 84
25) Isophorone	9.72	82	180575	48.07	ng	# 92
26) 2-Nitrophenol	9.92	139	43973	50.22	ng	# 77
27) 2,4-Dimethylphenol	9.98	122	83148	52.47	ng	89
28) bis(2-Chloroethoxy)methane	10.20	93	110164	51.19	ng	95
29) 2,4-Dichlorophenol	10.47	162	74295	50.61	ng	98
30) 1,2,4-Trichlorobenzene	10.67	180	69814	42.56	ng	95
31) Naphthalene	10.85	128	271335	48.56	ng	# 94
32) Benzoic acid	10.16	122	27238	49.33	ng	# 89
33) 4-Chloroaniline	10.99	127	26089	11.23	ng	# 87
34) Hexachlorobutadiene	11.12	225	34593	39.28	ng	97
35) Caprolactam	11.80	113	28356m	35.21	ng	
36) 4-Chloro-3-methylphenol	12.12	107	90528	52.02	ng	# 84
37) 2-Methylnaphthalene	12.46	142	195776	47.46	ng	95
39) 1,2,4,5-Tetrachlorobenzene	12.83	216	75371	42.32	ng	96
40) Hexachlorocyclopentadiene	12.80	237	29956	36.01	ng	96
42) 2,4,6-Trichlorophenol	13.08	196	57007	47.72	ng	98

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43) 2,4,5-Trichlorophenol	13.17	196	61597	46.67	nq	94
45) 1,1'-Biphenyl	13.46	154	253747	48.74	nq	94
46) 2-Chloronaphthalene	13.51	162	195593	49.01	nq	97
47) 2-Nitroaniline	13.73	65	50254	52.81	nq	# 58
48) Acenaphthylene	14.36	152	349694	49.94	nq	98
49) Dimethylphthalate	14.09	163	270523	47.16	nq	98
50) 2,6-Dinitrotoluene	14.22	165	60880	48.82	nq	# 81
51) Acenaphthene	14.70	154	207412	49.44	nq	94
52) 3-Nitroaniline	14.56	138	34923	25.25	nq	# 83
53) 2,4-Dinitrophenol	14.78	184	42804	83.50	nq	# 75
54) Dibenzofuran	15.03	168	319717	48.83	nq	98
55) 4-Nitrophenol	14.91	139	61971	64.92	nq	# 69
56) 2,4-Dinitrotoluene	15.01	165	84212	50.34	nq	# 90
57) Fluorene	15.68	166	271790	48.19	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.27	232	49313m	41.65	nq	
59) Diethylphthalate	15.44	149	294510	48.06	nq	98
60) 4-Chlorophenyl-phenylether	15.67	204	117179	45.88	nq	94
61) 4-Nitroaniline	15.73	138	55764m	38.01	nq	
62) Azobenzene	15.96	77	247884	53.42	nq	88
64) 4,6-Dinitro-2-methylphenol	15.78	198	36274	46.83	nq	# 84
65) n-Nitrosodiphenylamine	15.88	169	229242	51.92	nq	98
66) 4-Bromophenyl-phenylether	16.56	248	68087	47.41	nq	98
67) Hexachlorobenzene	16.68	284	69901	41.76	nq	98
68) Atrazine	16.83	200	69704	42.65	nq	96
69) Pentachlorophenol	17.04	266	41830	57.23	nq	96
70) Phenanthrene	17.42	178	414492	49.79	nq	97
71) Anthracene	17.52	178	410442	49.72	nq	97
72) Carbazole	17.79	167	378141	48.36	nq	96
73) Di-n-butylphthalate	18.32	149	523332	51.20	nq	98
74) Fluoranthene	19.43	202	449528	46.98	nq	94
77) Pyrene	19.79	202	449045	54.41	nq	98
79) Butylbenzylphthalate	20.66	149	235883	61.63	nq	# 91
80) Benzo(a)anthracene	21.62	228	377286	50.68	nq	99
81) 3,3'-Dichlorobenzidine	21.54	252	60485	28.94	nq	# 96
82) Chrysene	21.69	228	360945	49.83	nq	99
83) Bis(2-ethylhexyl)phthalate	21.49	149	335760	59.66	nq	# 95
84) Di-n-octyl phthalate	22.69	149	558117	59.72	nq	# 86
85) Indeno(1,2,3-cd)pyrene	28.52	276	375150	46.16	nq	# 78
87) Benzo(b)fluoranthene	23.82	252	349354	50.54	nq	# 94
88) Benzo(k)fluoranthene	23.89	252	347505	51.07	nq	# 92
89) Benzo(a)pyrene	24.69	252	338076	51.15	nq	# 94
90) Dibenzo(a,h)anthracene	28.55	278	304099	48.85	nq	# 87
91) Benzo(g,h,i)perylene	29.64	276	311406	48.08	nq	# 87

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

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