

Data Path : Z:\HPCHEM1\BNA G\DATA\BG022916\
 Data File : BG021155.D
 Acq On : 29 Feb 2016 14:55
 Operator : UM/SJ
 Sample : SSTDICCC040
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDICCC040

Quant Time: Feb 29 15:44:15 2016
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG022916.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 29 15:24:16 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.16	152	12147	20.00	ng	0.00
21) Naphthalene-d8	10.97	136	63287	20.00	ng	0.00
38) Acenaphthene-d10	14.76	164	40522	20.00	ng	0.00
63) Phenanthrene-d10	17.50	188	95053	20.00	ng	0.00
75) Chrysene-d12	21.77	240	103465	20.00	ng	0.00
86) Perylene-d12	25.04	264	93327	20.00	ng	0.00

System Monitoring Compounds						
5) 2-Fluorophenol	5.71	112	61592	88.77	ng	0.00
7) Phenol-d6	7.30	99	104138	100.86	ng	0.00
23) Nitrobenzene-d5	9.32	82	123502	92.98	ng	0.00
41) 2,4,6-Tribromophenol	16.24	330	34575	65.08	ng	0.00
44) 2-Fluorobiphenyl	13.39	172	233822	73.84	ng	0.00
78) Terphenyl-d14	20.09	244	350833	79.24	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	3.59	88	12892	40.13	ng	88
3) Pyridine	4.00	79	46092	49.05	ng	90
4) n-Nitrosodimethylamine	3.91	42	22998	48.27	ng	97
6) Aniline	7.48	93	68179	53.39	ng	# 88
8) 2-Chlorophenol	7.72	128	38407	44.38	ng	82
9) Benzaldehyde	7.29	77	29682	53.08	ng	83
10) Phenol	7.33	94	54308	51.35	ng	81
11) bis(2-Chloroethyl)ether	7.57	93	43293	53.40	ng	92
12) 1,3-Dichlorobenzene	8.05	146	38987	42.33	ng	# 90
13) 1,4-Dichlorobenzene	8.19	146	39717	40.37	ng	97
14) 1,2-Dichlorobenzene	8.51	146	39004	41.87	ng	97
15) Benzyl Alcohol	8.39	79	47973	53.19	ng	# 86
16) 2,2'-oxybis(1-Chloropropan	8.68	45	71883	77.33	ng	99
17) 2-Methylphenol	8.59	107	37895	50.76	ng	# 89
18) Hexachloroethane	9.25	117	16791	45.11	ng	96
19) n-Nitroso-di-n-propylamine	8.96	70	47265	59.35	ng	95
20) 3+4-Methylphenols	8.92	107	53336	51.53	ng	91
22) Acetophenone	8.98	105	76652	43.88	ng	# 92
24) Nitrobenzene	9.36	77	66324	48.33	ng	# 87
25) Isophorone	9.89	82	131641	52.88	ng	95
26) 2-Nitrophenol	10.07	139	24331	41.94	ng	# 54
27) 2,4-Dimethylphenol	10.12	122	43423	42.43	ng	92
28) bis(2-Chloroethoxy)methane	10.36	93	59955	49.42	ng	97
29) 2,4-Dichlorophenol	10.60	162	38620	38.55	ng	95
30) 1,2,4-Trichlorobenzene	10.83	180	39253	34.57	ng	93
31) Naphthalene	11.01	128	131944	40.59	ng	99
32) Benzoic acid	10.26	122	35277	37.54	ng	# 82
33) 4-Chloroaniline	11.12	127	61668	45.42	ng	# 89
34) Hexachlorobutadiene	11.30	225	24051	31.66	ng	89
35) Caprolactam	11.91	113	18946	55.97	ng	# 80
36) 4-Chloro-3-methylphenol	12.22	107	58140	47.34	ng	88
37) 2-Methylnaphthalene	12.61	142	93807	41.09	ng	# 92
39) 1,2,4,5-Tetrachlorobenzene	12.97	216	50582	33.78	ng	# 98
40) Hexachlorocyclopentadiene	12.95	237	28909	29.36	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.20	196	36839	39.23	nq	97
43) 2,4,5-Trichlorophenol	13.27	196	39636	41.08	nq	# 91
45) 1,1'-Biphenyl	13.61	154	135961	39.84	nq	97
46) 2-Chloronaphthalene	13.65	162	103167	39.85	nq	96
47) 2-Nitroaniline	13.85	65	49821	59.39	nq	# 74
48) Acenaphthylene	14.49	152	173714	42.40	nq	98
49) Dimethylphthalate	14.22	163	145548	41.68	nq	# 99
50) 2,6-Dinitrotoluene	14.33	165	30814	42.89	nq	# 58
51) Acenaphthene	14.83	154	112981	40.43	nq	98
52) 3-Nitroaniline	14.66	138	34123	48.81	nq	# 64
53) 2,4-Dinitrophenol	14.86	184	16618	33.17	nq	# 76
54) Dibenzofuran	15.16	168	144531	40.83	nq	92
55) 4-Nitrophenol	14.96	139	28889	46.59	nq	# 60
56) 2,4-Dinitrotoluene	15.12	165	44773	43.45	nq	# 71
57) Fluorene	15.81	166	139879	40.62	nq	97
58) 2,3,4,6-Tetrachlorophenol	15.38	232	31569	37.65	nq	# 92
59) Diethylphthalate	15.57	149	158997	43.19	nq	98
60) 4-Chlorophenyl-phenylether	15.80	204	69450	36.67	nq	97
61) 4-Nitroaniline	15.83	138	36832	46.27	nq	# 57
62) Azobenzene	16.08	77	173254	54.33	nq	92
64) 4,6-Dinitro-2-methylphenol	15.87	198	26645	38.22	nq	86
65) n-Nitrosodiphenylamine	16.01	169	120993	43.15	nq	94
66) 4-Bromophenyl-phenylether	16.69	248	41280	36.89	nq	# 89
67) Hexachlorobenzene	16.80	284	39779	33.69	nq	# 83
68) Atrazine	16.95	200	40883	39.34	nq	96
69) Pentachlorophenol	17.14	266	27096	34.00	nq	96
70) Phenanthrene	17.54	178	209365	40.94	nq	99
71) Anthracene	17.64	178	213801	41.51	nq	99
72) Carbazole	17.90	167	189508	42.62	nq	98
73) Di-n-butylphthalate	18.45	149	270405	46.13	nq	# 98
74) Fluoranthene	19.54	202	246777	38.46	nq	99
76) Benzidine	19.71	184	119613	43.94	nq	99
77) Pyrene	19.90	202	249148	43.49	nq	97
79) Butylbenzylphthalate	20.77	149	129788	54.28	nq	87
80) Benzo(a)anthracene	21.74	228	249895	41.98	nq	98
81) 3,3'-Dichlorobenzidine	21.65	252	95474	40.99	nq	98
82) Chrysene	21.81	228	236474	41.25	nq	98
83) Bis(2-ethylhexyl)phthalate	21.64	149	188890	51.16	nq	95
84) Di-n-octyl phthalate	22.87	149	316645	51.17	nq	99
85) Indeno(1,2,3-cd)pyrene	28.78	276	261206	37.79	nq	# 100
87) Benzo(b)fluoranthene	24.00	252	240017	41.14	nq	97
88) Benzo(k)fluoranthene	24.07	252	237870	41.37	nq	98
89) Benzo(a)pyrene	24.89	252	229700	41.02	nq	97
90) Dibenzo(a,h)anthracene	28.86	278	222874	39.48	nq	# 95
91) Benzo(g,h,i)perylene	29.96	276	212268	39.04	nq	# 92

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

