

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG062516\
 Data File : BG022578.D
 Acq On : 25 Jun 2016 17:26
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040EC

Quant Time: Jun 27 01:10:39 2016
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 24 20:14:59 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.17	152	133649	20.00	ng	0.00
21) Naphthalene-d8	10.99	136	572203	20.00	ng	0.00
38) Acenaphthene-d10	14.80	164	409075	20.00	ng	0.00
63) Phenanthrene-d10	17.56	188	1033603	20.00	ng	0.00
75) Chrysene-d12	21.86	240	1137831	20.00	ng	0.01
86) Perylene-d12	25.22	264	1201086	20.00	ng	0.01

System Monitoring Compounds

5) 2-Fluorophenol	5.70	112	639025	76.69	ng	0.00
7) Phenol-d6	7.31	99	949192	80.82	ng	0.00
23) Nitrobenzene-d5	9.34	82	1052921	77.88	ng	0.00
41) 2,4,6-Tribromophenol	16.29	330	535454	83.22	ng	0.00
44) 2-Fluorobiphenyl	13.42	172	2026158	78.55	ng	0.00
78) Terphenyl-d14	20.16	244	2974197	69.25	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.59	88	133998	39.68	ng	92
3) Pyridine	3.99	79	425373	45.03	ng	93
4) n-Nitrosodimethylamine	3.90	42	191516	35.44	ng	90
6) Aniline	7.48	93	683498	43.89	ng	95
8) 2-Chlorophenol	7.73	128	336149	38.95	ng	96
9) Benzaldehyde	7.30	77	182284	44.10	ng	92
10) Phenol	7.34	94	502014	40.44	ng	95
11) bis(2-Chloroethyl)ether	7.58	93	390249	38.19	ng	97
12) 1,3-Dichlorobenzene	8.05	146	397744	37.87	ng	96
13) 1,4-Dichlorobenzene	8.20	146	406717	38.63	ng	98
14) 1,2-Dichlorobenzene	8.52	146	380470	38.00	ng	97
15) Benzyl Alcohol	8.40	79	471849	43.42	ng	99
16) 2,2'-oxybis(1-Chloropropan	8.70	45	433309	38.34	ng	97
17) 2-Methylphenol	8.60	107	331579	40.52	ng	98
18) Hexachloroethane	9.26	117	164477	37.74	ng	88
19) n-Nitroso-di-n-propylamine	8.97	70	393449	40.22	ng	97
20) 3+4-Methylphenols	8.93	107	463461	41.12	ng	96
22) Acetophenone	8.99	105	646771	39.10	ng	# 94
24) Nitrobenzene	9.38	77	572019	38.28	ng	94
25) Isophorone	9.91	82	1026490	39.11	ng	100
26) 2-Nitrophenol	10.09	139	219098	40.71	ng	99
27) 2,4-Dimethylphenol	10.14	122	372131	36.19	ng	94
28) bis(2-Chloroethoxy)methane	10.38	93	548779	38.30	ng	96
29) 2,4-Dichlorophenol	10.63	162	413161	41.24	ng	94
30) 1,2,4-Trichlorobenzene	10.85	180	448747	37.72	ng	96
31) Naphthalene	11.04	128	1097796	38.17	ng	99
32) Benzoic acid	10.26	122	287418	43.32	ng	91
33) 4-Chloroaniline	11.15	127	525374	42.41	ng	96
34) Hexachlorobutadiene	11.32	225	355561	38.45	ng	95
35) Caprolactam	11.93	113	137799	43.66	ng	97
36) 4-Chloro-3-methylphenol	12.25	107	510596	41.52	ng	98
37) 2-Methylnaphthalene	12.64	142	881751	39.53	ng	97
39) 1,2,4,5-Tetrachlorobenzene	13.00	216	575577	37.27	ng	98
40) Hexachlorocyclopentadiene	12.98	237	450261	36.52	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	13.23	196	399319	40.32	ng	94
43) 2,4,5-Trichlorophenol	13.30	196	430217	41.52	ng	99
45) 1,1'-Biphenyl	13.64	154	1178661	37.82	ng	99
46) 2-Chloronaphthalene	13.68	162	977718	38.13	ng	97
47) 2-Nitroaniline	13.88	65	405736	40.99	ng	93
48) Acenaphthylene	14.52	152	1423541	38.28	ng	99
49) Dimethylphthalate	14.25	163	1300642	38.33	ng	97
50) 2,6-Dinitrotoluene	14.37	165	296783	40.25	ng	90
51) Acenaphthene	14.86	154	1087516	39.69	ng	97
52) 3-Nitroaniline	14.70	138	289675	41.98	ng	91
53) 2,4-Dinitrophenol	14.90	184	196983	43.37	ng	98
54) Dibenzofuran	15.20	168	1524165	38.41	ng	99
55) 4-Nitrophenol	14.99	139	250089	42.86	ng	96
56) 2,4-Dinitrotoluene	15.15	165	424863	41.59	ng	# 97
57) Fluorene	15.85	166	1255839	39.66	ng	97
58) 2,3,4,6-Tetrachlorophenol	15.42	232	451889	42.06	ng	97
59) Diethylphthalate	15.61	149	1348143	38.64	ng	97
60) 4-Chlorophenyl-phenylether	15.84	204	737436	39.12	ng	96
61) 4-Nitroaniline	15.86	138	298506	41.93	ng	89
62) Azobenzene	16.13	77	1447020	38.30	ng	93
64) 4,6-Dinitro-2-methylphenol	15.92	198	283716	40.93	ng	99
65) n-Nitrosodiphenylamine	16.05	169	1164382	38.71	ng	97
66) 4-Bromophenyl-phenylether	16.74	248	508391	37.91	ng	97
67) Hexachlorobenzene	16.86	284	545196	38.45	ng	97
68) Atrazine	17.00	200	511681	40.29	ng	99
69) Pentachlorophenol	17.20	266	399889	40.98	ng	95
70) Phenanthrene	17.60	178	2014981	38.23	ng	97
71) Anthracene	17.69	178	2062609	38.02	ng	98
72) Carbazole	17.96	167	1808793	39.33	ng	98
73) Di-n-butylphthalate	18.51	149	2112367	39.45	ng	99
74) Fluoranthene	19.61	202	2406664	39.63	ng	97
76) Benzidine	19.78	184	690427	56.99	ng	98
77) Pyrene	19.96	202	2376746	39.22	ng	96
79) Butylbenzylphthalate	20.85	149	1056799	40.28	ng	94
80) Benzo(a)anthracene	21.83	228	2540438	40.35	ng	99
81) 3,3'-Dichlorobenzidine	21.74	252	1012171	38.92	ng	97
82) Chrysene	21.90	228	2357206	39.84	ng	95
83) Bis(2-ethylhexyl)phthalate	21.73	149	1409678	38.16	ng	99
84) Di-n-octyl phthalate	22.99	149	2329248	38.25	ng	99
85) Indeno(1,2,3-cd)pyrene	29.07	276	3057406	39.32	ng	# 100
87) Benzo(b)fluoranthene	24.15	252	2742042	37.97	ng	98
88) Benzo(k)fluoranthene	24.21	252	2615796	38.31	ng	# 97
89) Benzo(a)pyrene	25.06	252	2667867	37.89	ng	99
90) Dibenzo(a,h)anthracene	29.15	278	2545231	38.40	ng	# 93
91) Benzo(g,h,i)perylene	30.28	276	2560060	37.94	ng	# 96

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

