

Data Path : Z:\SVOASRV\HPCHEM1\BNA G\DATA\BG060718\
 Data File : BG034951.D
 Acq On : 7 Jun 2018 13:56
 Operator : SJ/JU
 Sample : SSTDICV040
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 ICVBG060718

Quant Time: Jun 08 04:17:48 2018
 Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG060718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 07 13:13:20 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.08	152	35025	20.00	ng	0.00
21) Naphthalene-d8	10.89	136	143906	20.00	ng	0.00
38) Acenaphthene-d10	14.71	164	98228	20.00	ng	0.00
63) Phenanthrene-d10	17.45	188	256510	20.00	ng	0.00
75) Chrysene-d12	21.73	240	254188	20.00	ng	0.00
86) Perylene-d12	24.97	264	257117	20.00	ng	-0.01

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	5.65	112	168233	81.06	ng	0.00
7) Phenol-d6	7.23	99	254210	82.99	ng	0.00
23) Nitrobenzene-d5	9.24	82	262338	86.65	ng	0.00
41) 2,4,6-Tribromophenol	16.19	330	105566	83.95	ng	0.00
44) 2-Fluorobiphenyl	13.33	172	601364	79.59	ng	0.00
78) Terphenyl-d14	20.06	244	984602	81.14	ng	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.57	88	38680	39.062	ng	# 75
3) Pyridine	3.96	79	111004	41.547	ng	# 78
4) n-Nitrosodimethylamine	3.88	42	47149	41.077	ng	# 87
6) Aniline	7.41	93	158232	39.962	ng	99
8) 2-Chlorophenol	7.64	128	86451	39.739	ng	97
9) Benzaldehyde	7.22	77	95095	40.303	ng	98
10) Phenol	7.25	94	129935	40.988	ng	91
11) bis(2-Chloroethyl)ether	7.50	93	98416	39.999	ng	94
12) 1,3-Dichlorobenzene	7.97	146	104425	40.230	ng	# 94
13) 1,4-Dichlorobenzene	8.12	146	105896	39.521	ng	98
14) 1,2-Dichlorobenzene	8.43	146	101892	40.484	ng	96
15) Benzyl Alcohol	8.32	79	118321	40.765	ng	89
16) 2,2'-oxybis(1-Chloropropan	8.62	45	82002	33.485	ng	78
17) 2-Methylphenol	8.51	107	87544	41.887	ng	# 91
18) Hexachloroethane	9.17	117	38237	39.860	ng	86
19) n-Nitroso-di-n-propylamine	8.89	70	105094	40.383	ng	89
20) 3+4-Methylphenols	8.84	107	121836	40.670	ng	91
22) Acetophenone	8.90	105	179343	40.232	ng	93
24) Nitrobenzene	9.29	77	131520	42.424	ng	94
25) Isophorone	9.81	82	254437	39.807	ng	98
26) 2-Nitrophenol	10.00	139	47264	44.231	ng	# 85
27) 2,4-Dimethylphenol	10.05	122	87197	38.822	ng	92
28) bis(2-Chloroethoxy)methane	10.30	93	139494	40.611	ng	95
29) 2,4-Dichlorophenol	10.53	162	98752	40.407	ng	94
30) 1,2,4-Trichlorobenzene	10.75	180	115391	39.375	ng	96
31) Naphthalene	10.94	128	296815	39.704	ng	98
32) Benzoic acid	10.17	122	66848	44.433	ng	90
33) 4-Chloroaniline	11.04	127	131673	40.564	ng	94
34) Hexachlorobutadiene	11.22	225	90958	39.909	ng	97
35) Caprolactam	11.81	113	36258	40.143	ng	# 76
36) 4-Chloro-3-methylphenol	12.15	107	125564	41.216	ng	91
37) 2-Methylnaphthalene	12.54	142	225919	39.860	ng	# 90
39) 1,2,4,5-Tetrachlorobenzene	12.90	216	150600	41.161	ng	99
40) Hexachlorocyclopentadiene	12.89	237	93974	40.958	ng	89

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42) 2,4,6-Trichlorophenol	13.13	196	91863	41.931	nq	98
43) 2,4,5-Trichlorophenol	13.20	196	101482	42.203	nq	96
45) 1,1'-Biphenyl	13.54	154	315608	39.958	nq	98
46) 2-Chloronaphthalene	13.59	162	240126	40.210	nq	99
47) 2-Nitroaniline	13.78	65	80259	45.513	nq	97
48) Acenaphthylene	14.43	152	403003	40.246	nq	98
49) Dimethylphthalate	14.16	163	346093	40.404	nq	100
50) 2,6-Dinitrotoluene	14.27	165	68190	43.636	nq	91
51) Acenaphthene	14.77	154	271364	41.477	nq	100
52) 3-Nitroaniline	14.60	138	69646	45.419	nq	# 91
53) 2,4-Dinitrophenol	14.80	184	28574	45.181	nq	# 62
54) Dibenzofuran	15.11	168	394949	40.307	nq	96
55) 4-Nitrophenol	14.89	139	55984	43.260	nq	# 81
56) 2,4-Dinitrotoluene	15.05	165	93605	45.092	nq	# 95
57) Fluorene	15.75	166	332358	40.586	nq	98
58) 2,3,4,6-Tetrachlorophenol	15.32	232	98506	42.169	nq	98
59) Diethylphthalate	15.52	149	352868	40.377	nq	99
60) 4-Chlorophenyl-phenylether	15.75	204	188426	40.352	nq	97
61) 4-Nitroaniline	15.77	138	73793	44.873	nq	# 81
62) Azobenzene	16.04	77	347240	40.547	nq	96
64) 4,6-Dinitro-2-methylphenol	15.82	198	50141	42.813	nq	90
65) n-Nitrosodiphenylamine	15.96	169	307536	39.923	nq	99
66) 4-Bromophenyl-phenylether	16.64	248	121131	39.214	nq	92
67) Hexachlorobenzene	16.75	284	125201	39.821	nq	98
68) Atrazine	16.90	200	130320	39.671	nq	95
69) Pentachlorophenol	17.09	266	90325	42.723	nq	96
70) Phenanthrene	17.49	178	517165	39.690	nq	98
71) Anthracene	17.59	178	517972	39.685	nq	98
72) Carbazole	17.84	167	478599	39.520	nq	98
73) Di-n-butylphthalate	18.42	149	572686	39.675	nq	# 99
74) Fluoranthene	19.50	202	679395	40.069	nq	97
76) Benzidine	19.67	184	370663	39.196	nq	# 96
77) Pyrene	19.86	202	676560	40.374	nq	99
79) Butylbenzylphthalate	20.75	149	251822	41.385	nq	96
80) Benzo(a)anthracene	21.70	228	650748	40.062	nq	100
81) 3,3'-Dichlorobenzidine	21.62	252	252268	41.280	nq	99
82) Chrysene	21.77	228	595856	40.065	nq	99
83) Bis(2-ethylhexyl)phthalate	21.64	149	345048	40.132	nq	99
84) Di-n-octyl phthalate	22.88	149	602848	40.870	nq	# 94
85) Indeno(1,2,3-cd)pyrene	28.69	276	714687	41.031	nq	# 86
87) Benzo(b)fluoranthene	23.95	252	635753	40.152	nq	# 96
88) Benzo(k)fluoranthene	24.02	252	628184	40.557	nq	# 97
89) Benzo(a)pyrene	24.82	252	601315	40.359	nq	# 97
90) Dibenzo(a,h)anthracene	28.75	278	587161	39.921	nq	# 97
91) Benzo(g,h,i)perylene	29.83	276	580093	40.301	nq	# 94

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

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