

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070717\
 Data File : BG027870.D
 Acq On : 7 Jul 2017 16:01
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_G
 ClientSampleId :
 SSTDCCC040

Quant Time: Jul 07 17:30:34 2017
 Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 28 19:26:37 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	38765	20.00	ng	-0.05
21) Naphthalene-d8	11.25	136	191524	20.00	ng	-0.05
38) Acenaphthene-d10	15.02	164	110540	20.00	ng	-0.04
63) Phenanthrene-d10	17.77	188	265095	20.00	ng	-0.04
75) Chrysene-d12	22.13	240	315827	20.00	ng	-0.05
86) Perylene-d12	25.70	264	282696	20.00	ng	-0.09

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
5) 2-Fluorophenol	6.01	112	207203	82.31	ng	-0.05
7) Phenol-d6	7.57	99	302373	78.63	ng	-0.04
23) Nitrobenzene-d5	9.60	82	240777	70.08	ng	-0.05
41) 2,4,6-Tribromophenol	16.52	330	138538	91.37	ng	-0.03
44) 2-Fluorobiphenyl	13.65	172	644904	83.32	ng	-0.04
78) Terphenyl-d14	20.35	244	1027632	76.28	ng	-0.03

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
2) 1,4-Dioxane	3.98	88	40188	42.879	ng	# 65
3) Pyridine	4.37	79	110668	38.347	ng	# 63
4) n-Nitrosodimethylamine	4.28	42	31372	22.167	ng	# 1
6) Aniline	7.76	93	147130	33.005	ng	# 87
8) 2-Chlorophenol	8.00	128	113472	40.338	ng	# 79
9) Benzaldehyde	7.57	77	79091	34.562	ng	# 77
10) Phenol	7.60	94	146554	38.518	ng	# 66
11) bis(2-Chloroethyl)ether	7.84	93	121320	41.978	ng	# 76
12) 1,3-Dichlorobenzene	8.32	146	117509	39.273	ng	# 87
13) 1,4-Dichlorobenzene	8.46	146	121933	39.328	ng	95
14) 1,2-Dichlorobenzene	8.79	146	119507	39.754	ng	91
15) Benzyl Alcohol	8.66	79	39850	14.979	ng	# 74
16) 2,2'-oxybis(1-Chloropropan	8.94	45	133439	21.647	ng	83
17) 2-Methylphenol	8.86	107	102893	38.224	ng	# 81
18) Hexachloroethane	9.52	117	40268	36.924	ng	# 83
19) n-Nitroso-di-n-propylamine	9.22	70	91175	32.587	ng	# 73
20) 3+4-Methylphenols	9.18	107	142186	36.629	ng	85
22) Acetophenone	9.25	105	173683	37.580	ng	# 75
24) Nitrobenzene	9.64	77	129019	35.916	ng	# 78
25) Isophorone	10.15	82	252531	33.718	ng	# 92
26) 2-Nitrophenol	10.35	139	64301	39.629	ng	# 62
27) 2,4-Dimethylphenol	10.40	122	117726	38.737	ng	# 83
28) bis(2-Chloroethoxy)methane	10.62	93	158106	37.236	ng	100
29) 2,4-Dichlorophenol	10.89	162	104375	39.189	ng	95
30) 1,2,4-Trichlorobenzene	11.11	180	107235	39.478	ng	99
31) Naphthalene	11.30	128	397784	39.055	ng	98
32) Benzoic acid	10.49	122	19399	17.378	ng	# 76
33) 4-Chloroaniline	11.41	127	141936	32.257	ng	# 81
34) Hexachlorobutadiene	11.58	225	55413	36.612	ng	97
35) Caprolactam	12.18	113	45884	34.252	ng	# 50
36) 4-Chloro-3-methylphenol	12.49	107	120181	32.068	ng	# 83
37) 2-Methylnaphthalene	12.87	142	288356	38.011	ng	# 90
39) 1,2,4,5-Tetrachlorobenzene	13.23	216	126273	43.471	ng	98
40) Hexachlorocyclopentadiene	13.21	237	56787	41.464	ng	98

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42) 2,4,6-Trichlorophenol	13.47	196	84346	40.204	ng	93
43) 2,4,5-Trichlorophenol	13.55	196	95381	40.124	ng #	90
45) 1,1'-Biphenyl	13.86	154	359495	40.575	ng	98
46) 2-Chloronaphthalene	13.91	162	267780	40.008	ng	98
47) 2-Nitroaniline	14.11	65	80767	30.147	ng #	59
48) Acenaphthylene	14.75	152	495552	40.519	ng	98
49) Dimethylphthalate	14.47	163	340417	36.100	ng	97
50) 2,6-Dinitrotoluene	14.60	165	79059	38.568	ng #	62
51) Acenaphthene	15.09	154	311587	39.750	ng	97
52) 3-Nitroaniline	14.93	138	87615	36.331	ng #	78
53) 2,4-Dinitrophenol	15.13	184	37588	38.405	ng #	81
54) Dibenzofuran	15.42	168	414238	39.965	ng #	88
55) 4-Nitrophenol	15.23	139	59184	32.150	ng #	42
56) 2,4-Dinitrotoluene	15.38	165	109641	38.347	ng #	77
57) Fluorene	16.07	166	405848	40.384	ng	98
58) 2,3,4,6-Tetrachlorophenol	15.65	232	74465	36.339	ng #	96
59) Diethylphthalate	15.82	149	364050	34.865	ng	96
60) 4-Chlorophenyl-phenylether	16.05	204	175277	39.399	ng	93
61) 4-Nitroaniline	16.09	138	101389	40.356	ng #	49
62) Azobenzene	16.35	77	313630	34.182	ng	81
64) 4,6-Dinitro-2-methylphenol	16.14	198	59976	37.726	ng	83
65) n-Nitrosodiphenylamine	16.27	169	332212	38.524	ng	99
66) 4-Bromophenyl-phenylether	16.95	248	110454	39.837	ng	97
67) Hexachlorobenzene	17.08	284	128835	43.876	ng #	88
68) Atrazine	17.21	200	89434	32.849	ng	95
69) Pentachlorophenol	17.42	266	55498	33.489	ng	99
70) Phenanthrene	17.82	178	614445	40.149	ng	95
71) Anthracene	17.91	178	621050	40.187	ng	98
72) Carbazole	18.18	167	631636	41.501	ng #	94
73) Di-n-butylphthalate	18.70	149	658632	39.380	ng #	92
74) Fluoranthene	19.81	202	701938	41.968	ng	91
76) Benzidine	19.99	184	239490	32.938	ng	97
77) Pyrene	20.18	202	719847	36.319	ng	96
79) Butylbenzylphthalate	21.05	149	300881	33.706	ng	86
80) Benzo(a)anthracene	22.10	228	705049	36.924	ng	95
81) 3,3'-Dichlorobenzidine	22.00	252	268693	38.771	ng #	95
82) Chrysene	22.18	228	669342	37.398	ng	95
83) Bis(2-ethylhexyl)phthalate	21.96	149	448178	34.356	ng #	97
84) Di-n-octyl phthalate	23.29	149	753169	34.947	ng #	82
85) Indeno(1,2,3-cd)pyrene	29.80	276	825461	40.354	ng #	86
87) Benzo(b)fluoranthene	24.55	252	713231	39.161	ng #	94
88) Benzo(k)fluoranthene	24.63	252	726549	40.357	ng #	91
89) Benzo(a)pyrene	25.53	252	678295	39.690	ng #	93
90) Dibenzo(a,h)anthracene	29.87	278	719245	41.543	ng #	91
91) Benzo(g,h,i)perylene	31.09	276	673089	40.453	ng #	85

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

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