

Data Path : Z:\HPCHEM1\BNA F\DATA\BF041218\
 Data File : BF104446.D
 Acq On : 12 Apr 2018 13:02
 Operator : JU/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Quant Time: Apr 12 17:07:10 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 12 17:07:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.81	152	124516	20.00	ng	0.00
21) Naphthalene-d8	8.10	136	506895	20.00	ng	0.00
38) Acenaphthene-d10	9.85	164	240307	20.00	ng	0.00
63) Phenanthrene-d10	11.34	188	454356	20.00	ng	0.00
75) Chrysene-d12	13.99	240	370040	20.00	ng	0.00
86) Perylene-d12	15.46	264	317700	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.42	112	596208	73.21	ng	0.00
7) Phenol-d6	6.45	99	728635	72.82	ng	0.00
23) Nitrobenzene-d5	7.38	82	653602	84.20	ng	0.00
41) 2,4,6-Tribromophenol	10.64	330	203357	81.58	ng	0.00
44) 2-Fluorobiphenyl	9.17	172	1145682	75.32	ng	0.00
78) Terphenyl-d14	12.93	244	1195620	73.66	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	139662	36.807	ng	90
3) Pyridine	3.26	79	391025	36.653	ng	87
4) n-Nitrosodimethylamine	3.22	42	130067	34.878	ng	# 81
6) Aniline	6.47	93	506110	36.408	ng	98
8) 2-Chlorophenol	6.60	128	324959	37.808	ng	92
9) Benzaldehyde	6.36	77	184562	34.986	ng	96
10) Phenol	6.46	94	419325	39.499	ng	99
11) bis(2-Chloroethyl)ether	6.55	93	308398	36.403	ng	94
12) 1,3-Dichlorobenzene	6.75	146	369849	37.983	ng	100
13) 1,4-Dichlorobenzene	6.83	146	372427	38.369	ng	99
14) 1,2-Dichlorobenzene	6.98	146	342617	38.090	ng	99
15) Benzyl Alcohol	6.95	79	279149	36.733	ng	98
16) 2,2'-oxybis(1-Chloropropan	7.09	45	405218	35.617	ng	92
17) 2-Methylphenol	7.06	107	282254	37.779	ng	98
18) Hexachloroethane	7.32	117	133924	38.698	ng	96
19) n-Nitroso-di-n-propylamine	7.23	70	222527	34.035	ng	94
20) 3+4-Methylphenols	7.22	107	336903	36.359	ng	# 83
22) Acetophenone	7.22	105	452298	36.243	ng	95
24) Nitrobenzene	7.40	77	348584	40.672	ng	97
25) Isophorone	7.63	82	597868	36.421	ng	98
26) 2-Nitrophenol	7.71	139	175760	50.374	ng	97
27) 2,4-Dimethylphenol	7.75	122	254930	35.188	ng	99
28) bis(2-Chloroethoxy)methane	7.85	93	371438	37.008	ng	99
29) 2,4-Dichlorophenol	7.95	162	264604	38.279	ng	99
30) 1,2,4-Trichlorobenzene	8.03	180	289732	38.192	ng	99
31) Naphthalene	8.12	128	943685	37.387	ng	99
32) Benzoic acid	7.87	122	215566	37.292	ng	99
33) 4-Chloroaniline	8.17	127	409392	37.859	ng	96
34) Hexachlorobutadiene	8.23	225	162404	39.084	ng	99
35) Caprolactam	8.55	113	91450	37.924	ng	# 83
36) 4-Chloro-3-methylphenol	8.65	107	285879	38.570	ng	99
37) 2-Methylnaphthalene	8.81	142	613392	37.569	ng	100
39) 1,2,4,5-Tetrachlorobenzene	8.97	216	268380	39.015	ng	99
40) Hexachlorocyclopentadiene	8.96	237	156903	40.742	ng	98

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42) 2,4,6-Trichlorophenol	9.09	196	185776	38.904	nq	98
43) 2,4,5-Trichlorophenol	9.13	196	209643	39.232	nq	95
45) 1,1'-Biphenyl	9.27	154	751657	37.234	nq	99
46) 2-Chloronaphthalene	9.30	162	603114	37.823	nq	100
47) 2-Nitroaniline	9.40	65	186910	47.399	nq	95
48) Acenaphthylene	9.72	152	937983	37.492	nq	100
49) Dimethylphthalate	9.57	163	736884	38.886	nq	99
50) 2,6-Dinitrotoluene	9.64	165	158977	45.338	nq	97
51) Acenaphthene	9.89	154	550086	36.037	nq	99
52) 3-Nitroaniline	9.81	138	186631	45.464	nq	96
53) 2,4-Dinitrophenol	9.92	184	74252	58.904	nq	# 18
54) Dibenzofuran	10.06	168	797501	37.490	nq	98
55) 4-Nitrophenol	9.96	139	140950	43.710	nq	100
56) 2,4-Dinitrotoluene	10.05	165	211246	45.928	nq	90
57) Fluorene	10.40	166	612944	39.586	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.17	232	157896	39.606	nq	95
59) Diethylphthalate	10.27	149	715108	37.891	nq	98
60) 4-Chlorophenyl-phenylether	10.39	204	280750	40.714	nq	98
61) 4-Nitroaniline	10.42	138	188163	46.879	nq	99
62) Azobenzene	10.56	77	659251	36.562	nq	94
64) 4,6-Dinitro-2-methylphenol	10.45	198	116289	51.856	nq	90
65) n-Nitrosodiphenylamine	10.52	169	574250	36.452	nq	99
66) 4-Bromophenyl-phenylether	10.89	248	183001	37.866	nq	# 90
67) Hexachlorobenzene	10.95	284	207540	38.165	nq	91
68) Atrazine	11.04	200	181959	38.265	nq	99
69) Pentachlorophenol	11.14	266	129237	38.415	nq	97
70) Phenanthrene	11.37	178	926374	36.612	nq	99
71) Anthracene	11.42	178	930557	36.179	nq	99
72) Carbazole	11.57	167	914410	36.382	nq	99
73) Di-n-butylphthalate	11.90	149	1144924	37.292	nq	99
74) Fluoranthene	12.56	202	983789	37.213	nq	98
76) Benzidine	12.68	184	458295	35.312	nq	99
77) Pyrene	12.79	202	1008848	36.781	nq	100
79) Butylbenzylphthalate	13.40	149	497679	38.514	nq	99
80) Benzo(a)anthracene	13.98	228	846892	38.309	nq	100
81) 3,3'-Dichlorobenzidine	13.94	252	314171	38.116	nq	98
82) Chrysene	14.02	228	831554	36.621	nq	99
83) Bis(2-ethylhexyl)phthalate	13.96	149	668182	40.201	nq	99
84) Di-n-octyl phthalate	14.59	149	1028173	37.022	nq	95
85) Indeno(1,2,3-cd)pyrene	16.92	276	666418	34.549	nq	97
87) Benzo(b)fluoranthene	15.03	252	795962	39.668	nq	97
88) Benzo(k)fluoranthene	15.07	252	687447	36.289	nq	98
89) Benzo(a)pyrene	15.40	252	682465	38.013	nq	98
90) Dibenzo(a,h)anthracene	16.93	278	551234	37.384	nq	97
91) Benzo(g,h,i)perylene	17.36	276	537992	37.660	nq	94

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

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