

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040618\
 Data File : BF104300.D
 Acq On : 6 Apr 2018 12:27
 Operator : JU/SJ
 Sample : SSTDICC025
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC025

Quant Time: Apr 06 13:49:43 2018
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 06 13:46:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.80	152	137961	20.00	ng	0.00
21) Naphthalene-d8	8.09	136	560176	20.00	ng	0.00
38) Acenaphthene-d10	9.84	164	266709	20.00	ng	0.00
63) Phenanthrene-d10	11.33	188	490539	20.00	ng	0.00
75) Chrysene-d12	13.97	240	410571	20.00	ng	0.00
86) Perylene-d12	15.43	264	386016	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.40	112	445710	51.52	ng	0.00
7) Phenol-d6	6.42	99	552056	51.87	ng	0.00
23) Nitrobenzene-d5	7.37	82	420778	45.94	ng	0.00
41) 2,4,6-Tribromophenol	10.63	330	142854	48.10	ng	0.00
44) 2-Fluorobiphenyl	9.16	172	878926	51.97	ng	0.00
78) Terphenyl-d14	12.92	244	895536	50.36	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.52	88	102914	27.583	ng	91
3) Pyridine	3.24	79	283556	30.020	ng	88
4) n-Nitrosodimethylamine	3.19	42	99089	35.293	ng	# 87
6) Aniline	6.46	93	381500	29.942	ng	99
8) 2-Chlorophenol	6.58	128	241715	25.471	ng	97
9) Benzaldehyde	6.35	77	160272	28.372	ng	97
10) Phenol	6.43	94	293221	24.864	ng	98
11) bis(2-Chloroethyl)ether	6.54	93	234437	23.070	ng	95
12) 1,3-Dichlorobenzene	6.74	146	269557	25.270	ng	99
13) 1,4-Dichlorobenzene	6.82	146	269919	23.615	ng	97
14) 1,2-Dichlorobenzene	6.97	146	252489	24.577	ng	98
15) Benzyl Alcohol	6.94	79	213560	30.861	ng	97
16) 2,2'-oxybis(1-Chloropropan	7.08	45	312672	28.242	ng	95
17) 2-Methylphenol	7.05	107	206066	26.680	ng	98
18) Hexachloroethane	7.32	117	95763	25.024	ng	96
19) n-Nitroso-di-n-propylamine	7.22	70	175039	27.712	ng	96
20) 3+4-Methylphenols	7.20	107	263338	26.715	ng	# 78
22) Acetophenone	7.21	105	353647	26.093	ng	98
24) Nitrobenzene	7.39	77	237065	24.598	ng	97
25) Isophorone	7.62	82	444677	26.342	ng	98
26) 2-Nitrophenol	7.70	139	91004	18.089	ng	97
27) 2,4-Dimethylphenol	7.73	122	192997	26.600	ng	99
28) bis(2-Chloroethoxy)methane	7.83	93	278494	26.323	ng	99
29) 2,4-Dichlorophenol	7.93	162	192429	25.392	ng	99
30) 1,2,4-Trichlorobenzene	8.03	180	213368	24.815	ng	98
31) Naphthalene	8.11	128	715524	26.146	ng	100
32) Benzoic acid	7.83	122	142908	26.887	ng	98
33) 4-Chloroaniline	8.16	127	307207	26.627	ng	96
34) Hexachlorobutadiene	8.22	225	118904	25.412	ng	97
35) Caprolactam	8.52	113	67067	27.910	ng	# 85
36) 4-Chloro-3-methylphenol	8.63	107	207775	25.921	ng	99
37) 2-Methylnaphthalene	8.80	142	464382	25.761	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.96	216	196780	24.060	ng	100
40) Hexachlorocyclopentadiene	8.95	237	104843	31.117	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.07	196	131980	25.484	ng	98
43) 2,4,5-Trichlorophenol	9.10	196	152549	24.396	ng	97
45) 1,1'-Biphenyl	9.26	154	573520	25.054	ng	99
46) 2-Chloronaphthalene	9.29	162	451946	25.029	ng	99
47) 2-Nitroaniline	9.38	65	105840	20.556	ng	95
48) Acenaphthylene	9.70	152	716350	26.187	ng	99
49) Dimethylphthalate	9.56	163	531334	25.232	ng	100
50) 2,6-Dinitrotoluene	9.62	165	100381	22.157	ng	94
51) Acenaphthene	9.87	154	431164	27.246	ng	99
52) 3-Nitroaniline	9.79	138	116671	22.403	ng	# 96
53) 2,4-Dinitrophenol	9.89	184	21996	12.839	ng	# 23
54) Dibenzofuran	10.05	168	611467	26.460	ng	99
55) 4-Nitrophenol	9.94	139	91054	29.160	ng	94
56) 2,4-Dinitrotoluene	10.02	165	123314	21.331	ng	95
57) Fluorene	10.39	166	462419	24.258	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.16	232	113823	24.296	ng	96
59) Diethylphthalate	10.26	149	533888	26.466	ng	99
60) 4-Chlorophenyl-phenylether	10.38	204	206717	23.611	ng	97
61) 4-Nitroaniline	10.40	138	115811	23.766	ng	100
62) Azobenzene	10.55	77	513766	28.165	ng	94
64) 4,6-Dinitro-2-methylphenol	10.43	198	43237	14.561	ng	95
65) n-Nitrosodiphenylamine	10.50	169	434449	26.151	ng	99
66) 4-Bromophenyl-phenylether	10.87	248	131698	25.095	ng	92
67) Hexachlorobenzene	10.93	284	148958	24.675	ng	93
68) Atrazine	11.03	200	135217	26.568	ng	98
69) Pentachlorophenol	11.12	266	93699	28.206	ng	97
70) Phenanthrene	11.35	178	703826	26.501	ng	99
71) Anthracene	11.40	178	710930	25.627	ng	100
72) Carbazole	11.56	167	696628	26.311	ng	99
73) Di-n-butylphthalate	11.89	149	857569	27.192	ng	99
74) Fluoranthene	12.54	202	733555	25.985	ng	98
76) Benzidine	12.66	184	393327	33.617	ng	99
77) Pyrene	12.77	202	760552	25.769	ng	100
79) Butylbenzylphthalate	13.39	149	353863	25.449	ng	97
80) Benzo(a)anthracene	13.96	228	610643	23.770	ng	99
81) 3,3'-Dichlorobenzidine	13.93	252	235845	27.735	ng	96
82) Chrysene	14.00	228	635245	25.246	ng	98
83) Bis(2-ethylhexyl)phthalate	13.96	149	468205	26.060	ng	100
84) Di-n-octyl phthalate	14.57	149	780546	25.256	ng	97
85) Indeno(1,2,3-cd)pyrene	16.86	276	551271	21.142	ng	99
87) Benzo(b)fluoranthene	15.00	252	586268	24.955	ng	97
88) Benzo(k)fluoranthene	15.03	252	603114	27.253	ng	99
89) Benzo(a)pyrene	15.36	252	548604	25.200	ng	97
90) Dibenzo(a,h)anthracene	16.88	278	460387	22.873	ng	98
91) Benzo(g,h,i)perylene	17.29	276	437416	21.696	ng	98

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

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