

Data Path : Z:\HPCHEM1\BNA F\DATA\BF042815\
 Data File : BF078784.D
 Acq On : 28 Apr 2015 19:57
 Operator : TP/IZ
 Sample : SSTDICC060
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDICC060

Quant Time: Apr 29 03:21:44 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF042815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 29 03:08:11 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	90243	20.00	ng	0.00
21) Naphthalene-d8	8.97	136	377654	20.00	ng	0.00
38) Acenaphthene-d10	11.14	164	172982	20.00	ng	0.00
63) Phenanthrene-d10	12.99	188	334548	20.00	ng	0.00
75) Chrysene-d12	16.30	240	320964	20.00	ng	-0.01
86) Perylene-d12	18.08	264	320870	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.68	112	594795	115.68	ng	0.00
7) Phenol-d6	6.94	99	783448	121.96	ng	0.00
23) Nitrobenzene-d5	8.08	82	640509	111.67	ng	0.00
41) 2,4,6-Tribromophenol	12.12	330	193859	117.43	ng	0.00
44) 2-Fluorobiphenyl	10.32	172	1462399	122.52	ng	0.00
78) Terphenyl-d14	14.99	244	1531820	117.90	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.38	88	132193	56.14	ng	# 17
3) Pyridine	3.13	79	428492	68.08	ng	97
4) n-Nitrosodimethylamine	3.06	42	177212	64.66	ng	100
6) Aniline	6.98	93	579452	62.76	ng	97
8) 2-Chlorophenol	7.12	128	341650	55.91	ng	99
9) Benzaldehyde	6.84	77	284856	97.70	ng	97
10) Phenol	6.95	94	439136	57.34	ng	95
11) bis(2-Chloroethyl)ether	7.07	93	323731	57.09	ng	96
12) 1,3-Dichlorobenzene	7.31	146	377123	55.43	ng	99
13) 1,4-Dichlorobenzene	7.42	146	396761	56.15	ng	98
14) 1,2-Dichlorobenzene	7.60	146	374385	57.48	ng	99
15) Benzyl Alcohol	7.56	79	284389	56.85	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.75	45	514849	65.67	ng	99
17) 2-Methylphenol	7.70	107	276193	55.64	ng	98
18) Hexachloroethane	8.02	117	132534	56.81	ng	98
19) n-Nitroso-di-n-propylamine	7.91	70	281971	63.20	ng	# 87
20) 3+4-Methylphenols	7.90	107	359819	54.99	ng	94
22) Acetophenone	7.90	105	488534	58.00	ng	# 95
24) Nitrobenzene	8.10	77	333203	53.79	ng	97
25) Isophorone	8.40	82	688835	59.06	ng	99
26) 2-Nitrophenol	8.50	139	97361	34.78	ng	99
27) 2,4-Dimethylphenol	8.55	122	301034	54.78	ng	99
28) bis(2-Chloroethoxy)methane	8.68	93	429376	60.56	ng	100
29) 2,4-Dichlorophenol	8.79	162	274826	52.87	ng	98
30) 1,2,4-Trichlorobenzene	8.90	180	299820	51.90	ng	98
31) Naphthalene	8.99	128	998895	52.21	ng	100
32) Benzoic acid	8.65	122	100385	39.61	ng	98
33) 4-Chloroaniline	9.06	127	434471	55.72	ng	99
34) Hexachlorobutadiene	9.15	225	170672	51.93	ng	98
35) Caprolactam	9.50	113	84602	55.39	ng	92
36) 4-Chloro-3-methylphenol	9.66	107	283177	53.60	ng	97
37) 2-Methylnaphthalene	9.85	142	682967	53.19	ng	100
39) 1,2,4,5-Tetrachlorobenzene	10.06	216	278421	54.48	ng	# 100
40) Hexachlorocyclopentadiene	10.04	237	128571	59.57	ng	99

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	10.20	196	185583	53.70	nq	98
43) 2,4,5-Trichlorophenol	10.24	196	192028	50.91	nq	95
45) 1,1'-Biphenyl	10.43	154	766861	55.69	nq	99
46) 2-Chloronaphthalene	10.46	162	613840	54.92	nq	99
47) 2-Nitroaniline	10.58	65	150807	55.35	nq	95
48) Acenaphthylene	10.97	152	1017477	54.60	nq	99
49) Dimethylphthalate	10.82	163	722075	56.15	nq	98
50) 2,6-Dinitrotoluene	10.89	165	122540	47.55	nq	87
51) Acenaphthene	11.19	154	600432	56.39	nq	99
52) 3-Nitroaniline	11.10	138	168803	55.80	nq	89
53) 2,4-Dinitrophenol	11.22	184	18205	25.56	nq	87
54) Dibenzofuran	11.40	168	878922	55.61	nq	99
55) 4-Nitrophenol	11.29	139	124310	56.80	nq	93
56) 2,4-Dinitrotoluene	11.39	165	151112	45.60	nq	# 66
57) Fluorene	11.83	166	699910	53.57	nq	98
58) 2,3,4,6-Tetrachlorophenol	11.55	232	145055	50.11	nq	# 100
59) Diethylphthalate	11.70	149	714577	56.82	nq	98
60) 4-Chlorophenyl-phenylether	11.84	204	319149	53.81	nq	100
61) 4-Nitroaniline	11.85	138	157924	52.01	nq	93
62) Azobenzene	12.03	77	744989	62.61	nq	98
64) 4,6-Dinitro-2-methylphenol	11.88	198	34117	29.13	nq	79
65) n-Nitrosodiphenylamine	11.99	169	635426	57.74	nq	99
66) 4-Bromophenyl-phenylether	12.44	248	197663	55.99	nq	94
67) Hexachlorobenzene	12.50	284	215312	55.51	nq	97
68) Atrazine	12.65	200	170433	55.25	nq	98
69) Pentachlorophenol	12.75	266	100740	59.17	nq	99
70) Phenanthrene	13.03	178	1041957	54.81	nq	99
71) Anthracene	13.09	178	1010438	53.72	nq	100
72) Carbazole	13.29	167	988528	55.07	nq	99
73) Di-n-butylphthalate	13.74	149	1134425	56.43	nq	100
74) Fluoranthene	14.50	202	1036572	50.43	nq	100
76) Benzidine	14.69	184	662769	74.40	nq	99
77) Pyrene	14.79	202	1105971	56.25	nq	99
79) Butylbenzylphthalate	15.61	149	460430	58.74	nq	93
80) Benzo(a)anthracene	16.29	228	1026656	54.47	nq	100
81) 3,3'-Dichlorobenzidine	16.26	252	369571	59.12	nq	# 98
82) Chrysene	16.33	228	945647	55.70	nq	99
83) Bis(2-ethylhexyl)phthalate	16.34	149	718438	59.54	nq	100
84) Di-n-octyl phthalate	17.15	149	1119723	55.27	nq	# 100
85) Indeno(1,2,3-cd)pyrene	19.61	276	1233251	58.15	nq	# 100
87) Benzo(b)fluoranthene	17.62	252	988550	48.02	nq	# 97
88) Benzo(k)fluoranthene	17.66	252	1012985	61.28	nq	# 96
89) Benzo(a)pyrene	18.01	252	943454	54.50	nq	# 94
90) Dibenzo(a,h)anthracene	19.65	278	987728	57.00	nq	98
91) Benzo(g,h,i)perylene	20.07	276	980405	55.85	nq	97

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

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