

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051815\
 Data File : BF079338.D
 Acq On : 19 May 2015 7:48
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Quant Time: May 19 08:26:16 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 19 07:20:28 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.18	152	73806	20.00	ng	0.01
21) Naphthalene-d8	8.75	136	305589	20.00	ng	0.01
38) Acenaphthene-d10	10.93	164	151739	20.00	ng	0.01
63) Phenanthrene-d10	12.77	188	292840	20.00	ng	0.01
75) Chrysene-d12	16.06	240	293586	20.00	ng	0.00
86) Perylene-d12	17.82	264	306475	20.00	ng	-0.03

System Monitoring Compounds

5) 2-Fluorophenol	5.46	112	328210m	76.55	ng	0.01
7) Phenol-d6	6.74	99	435467	76.84	ng	0.00
23) Nitrobenzene-d5	7.87	82	417067	78.44	ng	0.01
41) 2,4,6-Tribromophenol	11.91	330	132957	81.00	ng	0.01
44) 2-Fluorobiphenyl	10.09	172	789251	72.47	ng	0.00
78) Terphenyl-d14	14.77	244	918794	74.93	ng	0.01

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.04	88	70700m	37.92	ng	
3) Pyridine	2.74	79	189164m	34.67	ng	
4) n-Nitrosodimethylamine	2.67	42	86756m	37.11	ng	
6) Aniline	6.76	93	298708	37.34	ng	# 40
8) 2-Chlorophenol	6.90	128	190576	39.19	ng	92
9) Benzaldehyde	6.62	77	130597	34.20	ng	97
10) Phenol	6.76	94	253909	38.62	ng	# 66
11) bis(2-Chloroethyl)ether	6.87	93	180255	37.40	ng	99
12) 1,3-Dichlorobenzene	7.10	146	210545	38.52	ng	97
13) 1,4-Dichlorobenzene	7.19	146	207072	36.81	ng	96
14) 1,2-Dichlorobenzene	7.37	146	192556	36.77	ng	96
15) Benzyl Alcohol	7.36	79	151291	35.96	ng	93
16) 2,2'-oxybis(1-Chloropropan	7.53	45	252654	34.26	ng	99
17) 2-Methylphenol	7.51	107	150411	38.43	ng	95
18) Hexachloroethane	7.79	117	60558	31.26	ng	95
19) n-Nitroso-di-n-propylamine	7.69	70	139948	36.56	ng	99
20) 3+4-Methylphenols	7.70	107	200708	38.49	ng	# 46
22) Acetophenone	7.68	105	258532	37.45	ng	# 91
24) Nitrobenzene	7.90	77	201041	38.14	ng	# 74
25) Isophorone	8.19	82	381496	38.70	ng	98
26) 2-Nitrophenol	8.28	139	88536	40.59	ng	97
27) 2,4-Dimethylphenol	8.35	122	178474	40.89	ng	98
28) bis(2-Chloroethoxy)methane	8.47	93	221994	36.53	ng	99
29) 2,4-Dichlorophenol	8.58	162	156719	40.37	ng	97
30) 1,2,4-Trichlorobenzene	8.68	180	166427	39.03	ng	98
31) Naphthalene	8.78	128	572148	39.29	ng	99
32) Benzoic acid	8.49	122	117920m	43.77	ng	
33) 4-Chloroaniline	8.86	127	251653	40.84	ng	# 89
34) Hexachlorobutadiene	8.92	225	89898	36.61	ng	99
35) Caprolactam	9.29	113	50444	40.19	ng	99
36) 4-Chloro-3-methylphenol	9.47	107	163378	40.07	ng	93
37) 2-Methylnaphthalene	9.63	142	387986	38.45	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.84	216	157923	36.95	ng	98
40) Hexachlorocyclopentadiene	9.83	237	5352	2.49	ng	97

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.99	196	111049	37.80	nq	97
43) 2,4,5-Trichlorophenol	10.04	196	112413	37.85	nq	98
45) 1,1'-Biphenyl	10.22	154	431925	35.81	nq	97
46) 2-Chloronaphthalene	10.24	162	356221	37.86	nq	93
47) 2-Nitroaniline	10.38	65	114666	42.06	nq	# 49
48) Acenaphthylene	10.74	152	586217	37.94	nq	99
49) Dimethylphthalate	10.60	163	396373	35.59	nq	99
50) 2,6-Dinitrotoluene	10.68	165	84115	38.27	nq	# 86
51) Acenaphthene	10.96	154	325694	35.35	nq	99
52) 3-Nitroaniline	10.88	138	116281	42.36	nq	95
53) 2,4-Dinitrophenol	11.03	184	1335	3.37	nq	# 75
54) Dibenzofuran	11.18	168	484428	36.40	nq	99
55) 4-Nitrophenol	11.11	139	76256	35.09	nq	# 87
56) 2,4-Dinitrotoluene	11.19	165	102901	35.42	nq	# 81
57) Fluorene	11.60	166	397674	36.41	nq	100
58) 2,3,4,6-Tetrachlorophenol	11.34	232	89554	36.89	nq	99
59) Diethylphthalate	11.49	149	405322	36.74	nq	98
60) 4-Chlorophenyl-phenylether	11.61	204	172819	35.67	nq	98
61) 4-Nitroaniline	11.65	138	131027	47.71	nq	85
62) Azobenzene	11.81	77	380934	35.04	nq	100
64) 4,6-Dinitro-2-methylphenol	11.69	198	4482	6.20	nq	80
65) n-Nitrosodiphenylamine	11.76	169	349990	35.89	nq	99
66) 4-Bromophenyl-phenylether	12.22	248	110804	36.30	nq	92
67) Hexachlorobenzene	12.29	284	126761	36.33	nq	# 66
68) Atrazine	12.43	200	82618	29.13	nq	97
69) Pentachlorophenol	12.54	266	62466	34.83	nq	97
70) Phenanthrene	12.79	178	587117	35.84	nq	99
71) Anthracene	12.86	178	586749	36.51	nq	100
72) Carbazole	13.06	167	561439	36.65	nq	99
73) Di-n-butylphthalate	13.52	149	665186	36.21	nq	100
74) Fluoranthene	14.27	202	647191	37.91	nq	94
76) Benzidine	14.46	184	426199	53.86	nq	97
77) Pyrene	14.55	202	651342	38.50	nq	99
79) Butylbenzylphthalate	15.37	149	301714	40.80	nq	88
80) Benzo(a)anthracene	16.05	228	621195	38.64	nq	99
81) 3,3'-Dichlorobenzidine	16.02	252	254312	44.41	nq	# 98
82) Chrysene	16.09	228	580010	37.51	nq	99
83) Bis(2-ethylhexyl)phthalate	16.10	149	420052	37.50	nq	99
84) Di-n-octyl phthalate	16.90	149	724715	36.74	nq	97
85) Indeno(1,2,3-cd)pyrene	19.22	276	739685	37.54	nq	100
87) Benzo(b)fluoranthene	17.36	252	677690	40.40	nq	100
88) Benzo(k)fluoranthene	17.39	252	548580	33.78	nq	# 95
89) Benzo(a)pyrene	17.75	252	577097	37.36	nq	# 98
90) Dibenzo(a,h)anthracene	19.25	278	609014	37.10	nq	97
91) Benzo(g,h,i)perylene	19.65	276	594871	36.16	nq	97

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

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