

Data Path : Z:\HPCHEM1\BNA F\DATA\BF052216\
 Data File : BF087299.D
 Acq On : 22 May 2016 19:13
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Quant Time: May 23 04:38:42 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF051716.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon May 23 03:47:59 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.56	152	145308	20.00	ng	0.00
21) Naphthalene-d8	7.85	136	611646	20.00	ng	0.00
38) Acenaphthene-d10	9.59	164	280955	20.00	ng	-0.01
63) Phenanthrene-d10	11.07	188	505943	20.00	ng	0.00
75) Chrysene-d12	13.70	240	304889	20.00	ng	0.00
86) Perylene-d12	15.05	264	275106	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.14	112	673435	77.90	ng	0.00
7) Phenol-d6	6.24	99	885610	76.38	ng	0.00
23) Nitrobenzene-d5	7.14	82	899349	73.28	ng	0.00
41) 2,4,6-Tribromophenol	10.39	330	209114	67.35	ng	0.00
44) 2-Fluorobiphenyl	8.92	172	1285714	75.79	ng	0.00
78) Terphenyl-d14	12.65	244	1230472	91.29	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.95	88	164167	36.92	ng	# 74
3) Pyridine	2.58	79	396005	36.81	ng	# 64
4) n-Nitrosodimethylamine	2.56	42	295548	38.77	ng	# 48
6) Aniline	6.23	93	536837	36.45	ng	98
8) 2-Chlorophenol	6.35	128	407670	38.76	ng	98
9) Benzaldehyde	6.11	77	225108	34.87	ng	97
10) Phenol	6.25	94	575512	39.49	ng	80
11) bis(2-Chloroethyl)ether	6.31	93	412964	37.73	ng	89
12) 1,3-Dichlorobenzene	6.49	146	428339m	38.31	ng	
13) 1,4-Dichlorobenzene	6.57	146	433356m	37.88	ng	
14) 1,2-Dichlorobenzene	6.73	146	393985	39.34	ng	99
15) Benzyl Alcohol	6.73	79	363587	42.12	ng	89
16) 2,2'-oxybis(1-Chloropropan	6.84	45	855074	38.01	ng	86
17) 2-Methylphenol	6.86	107	333637	38.62	ng	# 80
18) Hexachloroethane	7.06	117	174061	39.80	ng	96
19) n-Nitroso-di-n-propylamine	7.00	70	333793	37.87	ng	# 84
20) 3+4-Methylphenols	7.02	107	432987	38.00	ng	# 61
22) Acetophenone	6.99	105	632361	42.25	ng	96
24) Nitrobenzene	7.16	77	501641	37.44	ng	90
25) Isophorone	7.40	82	885058	39.40	ng	# 97
26) 2-Nitrophenol	7.47	139	212139	38.22	ng	# 61
27) 2,4-Dimethylphenol	7.54	122	350211	36.97	ng	94
28) bis(2-Chloroethoxy)methane	7.62	93	490043	39.73	ng	99
29) 2,4-Dichlorophenol	7.72	162	322812	38.67	ng	89
30) 1,2,4-Trichlorobenzene	7.79	180	355926	38.42	ng	97
31) Naphthalene	7.87	128	1167723	39.07	ng	99
32) Benzoic acid	7.71	122	209470	37.10	ng	# 79
33) 4-Chloroaniline	7.94	127	432132	34.79	ng	98
34) Hexachlorobutadiene	7.99	225	226789	43.14	ng	98
35) Caprolactam	8.34	113	103586	37.27	ng	# 64
36) 4-Chloro-3-methylphenol	8.44	107	384816	38.22	ng	89
37) 2-Methylnaphthalene	8.66	142	682577	35.71	ng	95
39) 1,2,4,5-Tetrachlorobenzene	8.73	216	348238	42.44	ng	99
40) Hexachlorocyclopentadiene	8.71	237	67581	27.09	ng	94

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42) 2,4,6-Trichlorophenol	8.86	196	206896	38.88	nq	92
43) 2,4,5-Trichlorophenol	8.90	196	214865	38.87	nq	# 85
45) 1,1'-Biphenyl	9.03	154	1032016	44.30	nq	98
46) 2-Chloronaphthalene	9.05	162	733002	40.99	nq	97
47) 2-Nitroaniline	9.16	65	272437	39.75	nq	# 79
48) Acenaphthylene	9.45	152	1039804	38.08	nq	98
49) Dimethylphthalate	9.34	163	823788	38.44	nq	99
50) 2,6-Dinitrotoluene	9.40	165	158276	34.01	nq	# 64
51) Acenaphthene	9.62	154	631767	37.03	nq	98
52) 3-Nitroaniline	9.58	138	173343	34.41	nq	# 75
53) 2,4-Dinitrophenol	9.70	184	57788	25.13	nq	# 83
54) Dibenzofuran	9.80	168	859929	38.98	nq	100
55) 4-Nitrophenol	9.78	139	24641m	15.02	nq	
56) 2,4-Dinitrotoluene	9.82	165	241811	40.57	nq	# 78
57) Fluorene	10.14	166	652893	36.83	nq	100
58) 2,3,4,6-Tetrachlorophenol	9.93	232	158002	36.70	nq	98
59) Diethylphthalate	10.03	149	825633	40.35	nq	98
60) 4-Chlorophenyl-phenylether	10.14	204	358529	42.67	nq	93
61) 4-Nitroaniline	10.19	138	159163	32.00	nq	# 63
62) Azobenzene	10.30	77	975550	41.38	nq	87
64) 4,6-Dinitro-2-methylphenol	10.23	198	98417	29.54	nq	89
65) n-Nitrosodiphenylamine	10.26	169	629071	39.51	nq	99
66) 4-Bromophenyl-phenylether	10.62	248	205911	39.58	nq	# 86
67) Hexachlorobenzene	10.68	284	225043	37.31	nq	# 75
68) Atrazine	10.80	200	206933	39.84	nq	93
69) Pentachlorophenol	10.89	266	56297	24.84	nq	97
70) Phenanthrene	11.10	178	1086527	39.60	nq	100
71) Anthracene	11.14	178	1017429	36.66	nq	100
72) Carbazole	11.31	167	890346	35.12	nq	98
73) Di-n-butylphthalate	11.64	149	1298840	44.60	nq	99
74) Fluoranthene	12.27	202	968890	35.24	nq	96
76) Benzidine	12.41	184	289359	31.45	nq	98
77) Pyrene	12.50	202	1001355	45.46	nq	100
79) Butylbenzylphthalate	13.13	149	467483	50.28	nq	93
80) Benzo(a)anthracene	13.69	228	676971	38.40	nq	98
81) 3,3'-Dichlorobenzidine	13.66	252	446228	39.78	nq	99
82) Chrysene	13.73	228	706567	41.43	nq	100
83) Bis(2-ethylhexyl)phthalate	13.69	149	600709	53.08	nq	# 99
84) Di-n-octyl phthalate	14.30	149	931300	47.77	nq	# 84
85) Indeno(1,2,3-cd)pyrene	16.27	276	648189	43.30	nq	94
87) Benzo(b)fluoranthene	14.69	252	584552m	32.03	nq	
88) Benzo(k)fluoranthene	14.71	252	635602m	46.78	nq	
89) Benzo(a)pyrene	14.99	252	590660	38.72	nq	# 96
90) Dibenzo(a,h)anthracene	16.29	278	543724	42.33	nq	# 92
91) Benzo(g,h,i)perylene	16.64	276	530184	40.87	nq	94

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

