

Data Path : Z:\HPCHEM1\BNA F\DATA\BF080917\
 Data File : BF097548.D
 Acq On : 10 Aug 2017 8:24
 Operator : SJ/JU
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 8/10/2017 6:08:09 PM

Quant Time: Aug 10 14:03:34 2017
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF072517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 07 16:02:51 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.40	152	90365	20.00	ng	-0.01
21) Naphthalene-d8	7.68	136	372409	20.00	ng	-0.02
38) Acenaphthene-d10	9.42	164	176510	20.00	ng	-0.02
63) Phenanthrene-d10	10.90	188	249170	20.00	ng	-0.02
75) Chrysene-d12	13.52	240	134686	20.00	ng	-0.02
86) Perylene-d12	14.86	264	163508	20.00	ng	-0.01

System Monitoring Compounds

5) 2-Fluorophenol	4.99	112	565182	94.28	ng	-0.02
7) Phenol-d6	6.07	99	596298	87.14	ng	-0.01
23) Nitrobenzene-d5	6.97	82	492800	77.63	ng	-0.02
41) 2,4,6-Tribromophenol	10.22	330	99681	71.48	ng	-0.02
44) 2-Fluorobiphenyl	8.76	172	841728	80.93	ng	-0.01
78) Terphenyl-d14	12.48	244	555880	84.15	ng	-0.02

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.86	88	96575	38.607	ng	# 73
3) Pyridine	2.45	79	299336	39.078	ng	# 60
4) n-Nitrosodimethylamine	2.42	42	171703	40.462	ng	81
6) Aniline	6.06	93	392926	45.628	ng	97
8) 2-Chlorophenol	6.19	128	288891	45.071	ng	91
9) Benzaldehyde	5.95	77	223349	55.140	ng	96
10) Phenol	6.09	94	367002	45.213	ng	96
11) bis(2-Chloroethyl)ether	6.15	93	281300	43.816	ng	90
12) 1,3-Dichlorobenzene	6.33	146	298531	43.218	ng	97
13) 1,4-Dichlorobenzene	6.42	146	292126	42.674	ng	97
14) 1,2-Dichlorobenzene	6.57	146	246513	41.243	ng	98
15) Benzyl Alcohol	6.57	79	194843	44.970	ng	96
16) 2,2'-oxybis(1-Chloropropan	6.69	45	508176	39.465	ng	86
17) 2-Methylphenol	6.69	107	199617	40.352	ng	# 88
18) Hexachloroethane	6.90	117	97773	39.041	ng	# 80
19) n-Nitroso-di-n-propylamine	6.83	70	185965	41.284	ng	# 81
20) 3+4-Methylphenols	6.85	107	279846	43.313	ng	# 61
22) Acetophenone	6.82	105	355664	39.769	ng	# 99
24) Nitrobenzene	6.99	77	270132	40.858	ng	93
25) Isophorone	7.24	82	532875	42.863	ng	97
26) 2-Nitrophenol	7.31	139	159796	44.674	ng	# 86
27) 2,4-Dimethylphenol	7.37	122	250844	42.512	ng	98
28) bis(2-Chloroethoxy)methane	7.45	93	319218	39.347	ng	98
29) 2,4-Dichlorophenol	7.57	162	216221	42.537	ng	96
30) 1,2,4-Trichlorobenzene	7.63	180	192899	39.813	ng	98
31) Naphthalene	7.70	128	769650	43.842	ng	100
32) Benzoic acid	7.56	122	139378	31.634	ng	91
33) 4-Chloroaniline	7.77	127	326442	45.701	ng	99
34) Hexachlorobutadiene	7.83	225	111264	43.317	ng	97
35) Caprolactam	8.15	113	72590m	44.535	ng	
36) 4-Chloro-3-methylphenol	8.28	107	253423	45.698	ng	90
37) 2-Methylnaphthalene	8.39	142	483524	43.502	ng	99
39) 1,2,4,5-Tetrachlorobenzene	8.56	216	189298	40.193	ng	99
40) Hexachlorocyclopentadiene	8.55	237	12628	13.627	ng	92

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42) 2,4,6-Trichlorophenol	8.69	196	132786	38.906	nq	99
43) 2,4,5-Trichlorophenol	8.75	196	123598	36.180	nq	95
45) 1,1'-Biphenyl	8.86	154	626876	41.580	nq	98
46) 2-Chloronaphthalene	8.88	162	454122	40.932	nq	# 92
47) 2-Nitroaniline	8.99	65	178350	44.296	nq	97
48) Acenaphthylene	9.29	152	692931	40.691	nq	98
49) Dimethylphthalate	9.17	163	559638	41.752	nq	99
50) 2,6-Dinitrotoluene	9.24	165	114138	40.149	nq	# 79
51) Acenaphthene	9.46	154	411080	40.982	nq	100
52) 3-Nitroaniline	9.40	138	143206	40.583	nq	94
53) 2,4-Dinitrophenol	9.55	184	25797	19.958	nq	# 74
54) Dibenzofuran	9.63	168	566546	42.403	nq	93
55) 4-Nitrophenol	9.61	139	61192m	29.160	nq	
56) 2,4-Dinitrotoluene	9.65	165	150553	46.067	nq	# 68
57) Fluorene	9.97	166	414835	41.310	nq	98
58) 2,3,4,6-Tetrachlorophenol	9.77	232	87151	36.948	nq	# 97
59) Diethylphthalate	9.87	149	547925	44.557	nq	100
60) 4-Chlorophenyl-phenylether	9.97	204	179611	43.314	nq	96
61) 4-Nitroaniline	10.02	138	127715	38.091	nq	93
62) Azobenzene	10.13	77	492017	43.129	nq	92
64) 4,6-Dinitro-2-methylphenol	10.06	198	38624m	24.946	nq	
65) n-Nitrosodiphenylamine	10.09	169	403219	46.509	nq	97
66) 4-Bromophenyl-phenylether	10.45	248	114350	45.986	nq	96
67) Hexachlorobenzene	10.52	284	120074	43.060	nq	# 90
68) Atrazine	10.63	200	105869	42.585	nq	94
69) Pentachlorophenol	10.73	266	74890m	64.909	nq	
70) Phenanthrene	10.92	178	548838	41.110	nq	100
71) Anthracene	10.97	178	555127	40.597	nq	98
72) Carbazole	11.14	167	532241	39.578	nq	99
73) Di-n-butylphthalate	11.48	149	756804	45.527	nq	98
74) Fluoranthene	12.10	202	441764	33.777	nq	98
76) Benzidine	12.24	184	143164	28.647	nq	# 93
77) Pyrene	12.33	202	455138	41.277	nq	97
79) Butylbenzylphthalate	12.96	149	242917	43.310	nq	# 78
80) Benzo(a)anthracene	13.51	228	343558	39.423	nq	99
81) 3,3'-Dichlorobenzidine	13.48	252	133407	40.594	nq	# 97
82) Chrysene	13.55	228	381640	44.848	nq	100
83) Bis(2-ethylhexyl)phthalate	13.52	149	299173	40.853	nq	98
84) Di-n-octyl phthalate	14.13	149	616964	45.909	nq	# 91
85) Indeno(1,2,3-cd)pyrene	16.04	276	314836	43.147	nq	96
87) Benzo(b)fluoranthene	14.52	252	413746	42.095	nq	99
88) Benzo(k)fluoranthene	14.54	252	336644m	35.943	nq	
89) Benzo(a)pyrene	14.81	252	360246	40.047	nq	99
90) Dibenzo(a,h)anthracene	16.04	278	264066	35.797	nq	96
91) Benzo(g,h,i)perylene	16.39	276	257184	33.246	nq	98

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

