

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060315\
 Data File : BF079674.D
 Acq On : 3 Jun 2015 21:07
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

apatel
 6/5/2015 8:06:31 AM

Quant Time: Jun 05 11:39:26 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 05 10:52:02 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.52	152	85649	20.00	ng	0.00
21) Naphthalene-d8	8.81	136	329701	20.00	ng	0.00
38) Acenaphthene-d10	10.59	164	168852	20.00	ng	0.00
63) Phenanthrene-d10	12.11	188	352178	20.00	ng	-0.02
75) Chrysene-d12	15.19	240	372882	20.00	ng	-0.19
86) Perylene-d12	17.19	264	338985	20.00	ng	-0.30

System Monitoring Compounds

5) 2-Fluorophenol	6.14	112	380601	80.69	ng	0.00
7) Phenol-d6	7.11	99	533734	86.06	ng	0.00
23) Nitrobenzene-d5	8.08	82	455886	81.07	ng	0.00
41) 2,4,6-Tribromophenol	11.38	330	115086	82.58	ng	-0.01
44) 2-Fluorobiphenyl	9.88	172	971425	87.28	ng	0.00
78) Terphenyl-d14	13.85	244	1267014	86.27	ng	-0.10

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.60	88	86935	39.80	ng	99
3) Pyridine	4.31	79	249876	39.80	ng	99
4) n-Nitrosodimethylamine	4.23	42	118300	42.27	ng	99
6) Aniline	7.18	93	364507	40.94	ng	99
8) 2-Chlorophenol	7.30	128	230835	42.04	ng	97
9) Benzaldehyde	7.06	77	149422	41.55	ng	99
10) Phenol	7.12	94	281821	42.37	ng	99
11) bis(2-Chloroethyl)ether	7.23	93	218490	41.20	ng	97
12) 1,3-Dichlorobenzene	7.46	146	256688	40.88	ng	99
13) 1,4-Dichlorobenzene	7.54	146	259646	40.44	ng	98
14) 1,2-Dichlorobenzene	7.69	146	237776m	40.91	ng	
15) Benzyl Alcohol	7.63	79	182159	40.28	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.78	45	348098	39.37	ng	99
17) 2-Methylphenol	7.74	107	184039	40.60	ng	98
18) Hexachloroethane	8.04	117	83932	39.93	ng	93
19) n-Nitroso-di-n-propylamine	7.91	70	179967	42.39	ng	98
20) 3+4-Methylphenols	7.88	107	245136	40.83	ng	99
22) Acetophenone	7.92	105	313544	40.64	ng	97
24) Nitrobenzene	8.09	77	229669	41.57	ng	98
25) Isophorone	8.33	82	467717	41.62	ng	99
26) 2-Nitrophenol	8.41	139	73554	40.31	ng	96
27) 2,4-Dimethylphenol	8.43	122	186999m	36.36	ng	
28) bis(2-Chloroethoxy)methane	8.52	93	259306	40.16	ng	99
29) 2,4-Dichlorophenol	8.65	162	195807	42.69	ng	97
30) 1,2,4-Trichlorobenzene	8.75	180	205027	39.66	ng	98
31) Naphthalene	8.83	128	685647	39.78	ng	100
32) Benzoic acid	8.48	122	65484m	45.49	ng	
33) 4-Chloroaniline	8.87	127	385493	41.13	ng	98
34) Hexachlorobutadiene	8.95	225	113112	39.04	ng	95
35) Caprolactam	9.20	113	56170	44.48	ng	# 62
36) 4-Chloro-3-methylphenol	9.32	107	198175	42.33	ng	97
37) 2-Methylnaphthalene	9.53	142	478052	41.58	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.70	216	224356	41.40	ng	98
40) Hexachlorocyclopentadiene	9.69	237	93351	36.13	ng	100

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42) 2,4,6-Trichlorophenol	9.80	196	126080	42.79	nq	99
43) 2,4,5-Trichlorophenol	9.84	196	150406	46.07	nq	96
45) 1,1'-Biphenyl	9.99	154	534388	41.04	nq	95
46) 2-Chloronaphthalene	10.02	162	426837	41.84	nq	# 90
47) 2-Nitroaniline	10.10	65	101561	48.35	nq	92
48) Acenaphthylene	10.46	152	702638	41.55	nq	99
49) Dimethylphthalate	10.27	163	513213	41.48	nq	99
50) 2,6-Dinitrotoluene	10.34	165	92002	46.81	nq	91
51) Acenaphthene	10.63	154	426226	41.59	nq	98
52) 3-Nitroaniline	10.51	138	120083	49.53	nq	93
53) 2,4-Dinitrophenol	10.62	184	22583	40.41	nq	# 1
54) Dibenzofuran	10.80	168	629602	42.33	nq	97
55) 4-Nitrophenol	10.65	139	82677	42.28	nq	94
56) 2,4-Dinitrotoluene	10.75	165	105079	43.57	nq	88
57) Fluorene	11.14	166	527213	43.37	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.90	232	107326	47.44	nq	99
59) Diethylphthalate	10.98	149	538538	45.33	nq	97
60) 4-Chlorophenyl-phenylether	11.13	204	235603	42.94	nq	91
61) 4-Nitroaniline	11.13	138	118629	49.60	nq	95
62) Azobenzene	11.29	77	466779	42.02	nq	86
64) 4,6-Dinitro-2-methylphenol	11.16	198	38469	39.34	nq	# 73
65) n-Nitrosodiphenylamine	11.23	169	440245	39.81	nq	99
66) 4-Bromophenyl-phenylether	11.62	248	145455	41.16	nq	95
67) Hexachlorobenzene	11.71	284	164257	39.56	nq	95
68) Atrazine	11.76	200	127953	45.17	nq	96
69) Pentachlorophenol	11.90	266	72912	47.93	nq	98
70) Phenanthrene	12.14	178	812021	41.70	nq	100
71) Anthracene	12.19	178	798261	40.75	nq	99
72) Carbazole	12.34	167	743644	41.34	nq	98
73) Di-n-butylphthalate	12.67	149	847560	42.99	nq	# 79
74) Fluoranthene	13.44	202	865111	41.21	nq	93
76) Benzidine	13.55	184	297645	35.09	nq	99
77) Pyrene	13.70	202	880185	41.00	nq	99
79) Butylbenzylphthalate	14.41	149	329070	45.02	nq	97
80) Benzo(a)anthracene	15.17	228	857445	42.05	nq	99
81) 3,3'-Dichlorobenzidine	15.11	252	289406	47.51	nq	98
82) Chrysene	15.22	228	843601	41.12	nq	98
83) Bis(2-ethylhexyl)phthalate	15.13	149	528158	43.49	nq	# 97
84) Di-n-octyl phthalate	15.98	149	816730	45.58	nq	100
85) Indeno(1,2,3-cd)pyrene	19.22	276	909447	45.33	nq	100
87) Benzo(b)fluoranthene	16.59	252	790908	45.15	nq	98
88) Benzo(k)fluoranthene	16.64	252	846301	41.24	nq	99
89) Benzo(a)pyrene	17.11	252	768313	44.86	nq	# 97
90) Dibenzo(a,h)anthracene	19.26	278	750618	46.77	nq	99
91) Benzo(g,h,i)perylene	19.84	276	758081	44.53	nq	98

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

