

Data Path : Z:\HPCHEM1\BNA F\DATA\BF060515\
 Data File : BF079767.D
 Acq On : 6 Jun 2015 12:42
 Operator : TP/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

apatel
 6/9/2015 8:16:21 AM

Quant Time: Jun 08 17:29:32 2015
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 08 15:29:07 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.51	152	46083	20.00	ng	0.01
21) Naphthalene-d8	8.79	136	195880	20.00	ng	0.00
38) Acenaphthene-d10	10.56	164	109939	20.00	ng	0.00
63) Phenanthrene-d10	12.11	188	204785	20.00	ng	0.05
75) Chrysene-d12	15.34	240	193308m	20.00	ng	0.40
86) Perylene-d12	17.44	264	179541m	20.00	ng	0.62

System Monitoring Compounds

5) 2-Fluorophenol	6.18	112	214349	77.96	ng	0.08
7) Phenol-d6	7.10	99	262625	73.95	ng	0.01
23) Nitrobenzene-d5	8.06	82	239408	75.04	ng	0.01
41) 2,4,6-Tribromophenol	11.37	330	89576	85.19	ng	0.01
44) 2-Fluorobiphenyl	9.86	172	547372	67.45	ng	0.00
78) Terphenyl-d14	13.93	244	729120m	86.38	ng	0.22

Target Compounds

						Qvalue
2) 1,4-Dioxane	4.20	88	65177m	53.57	ng	
3) Pyridine	4.72	79	127179m	37.37	ng	
4) n-Nitrosodimethylamine	4.66	42	60223m	40.78	ng	
6) Aniline	7.16	93	190626	36.99	ng	96
8) 2-Chlorophenol	7.29	128	110694	37.19	ng	91
9) Benzaldehyde	7.06	77	84956	36.94	ng	97
10) Phenol	7.12	94	135578	35.07	ng	99
11) bis(2-Chloroethyl)ether	7.22	93	117397	39.16	ng	97
12) 1,3-Dichlorobenzene	7.45	146	137316	36.85	ng	97
13) 1,4-Dichlorobenzene	7.53	146	145504	36.80	ng	98
14) 1,2-Dichlorobenzene	7.68	146	135510m	36.69	ng	
15) Benzyl Alcohol	7.62	79	109624	37.35	ng	95
16) 2,2'-oxybis(1-Chloropropan	7.76	45	193932	42.78	ng	98
17) 2-Methylphenol	7.72	107	100541	36.22	ng	98
18) Hexachloroethane	8.02	117	32488	28.28	ng	92
19) n-Nitroso-di-n-propylamine	7.88	70	112696	45.74	ng	96
20) 3+4-Methylphenols	7.87	107	145496	41.13	ng	99
22) Acetophenone	7.90	105	191146	40.91	ng	# 99
24) Nitrobenzene	8.07	77	116413	36.58	ng	98
25) Isophorone	8.31	82	253857	38.98	ng	96
26) 2-Nitrophenol	8.39	139	31736	23.68	ng	# 63
27) 2,4-Dimethylphenol	8.41	122	104068	34.50	ng	96
28) bis(2-Chloroethoxy)methane	8.50	93	155460	38.63	ng	99
29) 2,4-Dichlorophenol	8.63	162	106114	39.22	ng	97
30) 1,2,4-Trichlorobenzene	8.73	180	114459	34.83	ng	97
31) Naphthalene	8.81	128	401677	39.09	ng	99
32) Benzoic acid	8.46	122	48791	42.44	ng	92
33) 4-Chloroaniline	8.84	127	195250	37.40	ng	96
34) Hexachlorobutadiene	8.92	225	76636	38.34	ng	95
35) Caprolactam	9.18	113	31220	37.88	ng	99
36) 4-Chloro-3-methylphenol	9.30	107	104842	34.68	ng	# 77
37) 2-Methylnaphthalene	9.51	142	258317	39.54	ng	99
39) 1,2,4,5-Tetrachlorobenzene	9.68	216	116361	33.99	ng	99
40) Hexachlorocyclopentadiene	9.67	237	4919	3.34	ng	96

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42) 2,4,6-Trichlorophenol	9.78	196	74396	32.67	nq	98
43) 2,4,5-Trichlorophenol	9.82	196	88773	37.05	nq	96
45) 1,1'-Biphenyl	9.96	154	307314	32.13	nq	95
46) 2-Chloronaphthalene	10.00	162	253313	32.44	nq	97
47) 2-Nitroaniline	10.08	65	66058	40.33	nq	89
48) Acenaphthylene	10.42	152	441249	34.18	nq	99
49) Dimethylphthalate	10.25	163	299322	35.35	nq	# 98
50) 2,6-Dinitrotoluene	10.32	165	42306	32.75	nq	# 80
51) Acenaphthene	10.59	154	283042	38.71	nq	98
52) 3-Nitroaniline	10.49	138	76524	42.79	nq	92
53) 2,4-Dinitrophenol	10.59	184	817	5.37	nq	# 1
54) Dibenzofuran	10.77	168	417731	40.34	nq	91
55) 4-Nitrophenol	10.63	139	62035	46.43	nq	95
56) 2,4-Dinitrotoluene	10.73	165	56086	35.77	nq	# 94
57) Fluorene	11.12	166	383274	42.76	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.88	232	81669	42.11	nq	99
59) Diethylphthalate	10.96	149	352101	42.52	nq	96
60) 4-Chlorophenyl-phenylether	11.10	204	156069	38.42	nq	96
61) 4-Nitroaniline	11.11	138	55015	33.30	nq	96
62) Azobenzene	11.27	77	341030	40.75	nq	82
64) 4,6-Dinitro-2-methylphenol	11.14	198	1869	5.65	nq	# 30
65) n-Nitrosodiphenylamine	11.21	169	277218	40.53	nq	99
66) 4-Bromophenyl-phenylether	11.61	248	88062	39.71	nq	92
67) Hexachlorobenzene	11.70	284	98779	39.61	nq	# 96
68) Atrazine	11.75	200	82575	40.17	nq	98
69) Pentachlorophenol	11.90	266	45276	42.69	nq	98
70) Phenanthrene	12.14	178	492453	41.07	nq	98
71) Anthracene	12.19	178	491705	41.59	nq	99
72) Carbazole	12.34	167	437074	38.90	nq	98
73) Di-n-butylphthalate	12.69	149	509967	40.04	nq	# 82
74) Fluoranthene	13.49	202	538150	40.99	nq	98
76) Benzidine	13.61	184	17094m	2.80	nq	
77) Pyrene	13.77	202	557522m	47.04	nq	
79) Butylbenzylphthalate	14.53	149	216814m	45.89	nq	
80) Benzo(a)anthracene	15.33	228	497501m	42.04	nq	
81) 3,3'-Dichlorobenzidine	15.27	252	92705m	23.06	nq	
82) Chrysene	15.38	228	429940m	39.30	nq	
83) Bis(2-ethylhexyl)phthalate	15.30	149	317937m	42.59	nq	
84) Di-n-octyl phthalate	16.22	149	539993m	41.07	nq	
85) Indeno(1,2,3-cd)pyrene	19.47	276	432511m	34.16	nq	
87) Benzo(b)fluoranthene	16.83	252	445880m	39.98	nq	
88) Benzo(k)fluoranthene	16.88	252	434565m	40.62	nq	
89) Benzo(a)pyrene	17.35	252	411132m	39.93	nq	
90) Dibenzo(a,h)anthracene	19.51	278	352469m	35.47	nq	
91) Benzo(g,h,i)perylene	20.09	276	349363m	34.01	nq	

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

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