

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040816\
 Data File : BF086163.D
 Acq On : 8 Apr 2016 12:31
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

Sohil
 4/11/2016 5:45:06 PM

Quant Time: Apr 08 23:30:19 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 01 23:17:06 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.73	152	141260	20.00	ng	-0.05
21) Naphthalene-d8	8.02	136	594975	20.00	ng	-0.05
38) Acenaphthene-d10	9.77	164	266421	20.00	ng	-0.05
63) Phenanthrene-d10	11.25	188	515000	20.00	ng	-0.05
75) Chrysene-d12	13.89	240	401320	20.00	ng	-0.03
86) Perylene-d12	15.28	264	307809	20.00	ng	-0.06

System Monitoring Compounds

5) 2-Fluorophenol	5.32	112	713574	86.35	ng	-0.05
7) Phenol-d6	6.38	99	799300	76.14	ng	-0.03
23) Nitrobenzene-d5	7.31	82	853189	84.57	ng	-0.03
41) 2,4,6-Tribromophenol	10.57	330	206368	82.53	ng	-0.03
44) 2-Fluorobiphenyl	9.09	172	1183043	73.83	ng	-0.05
78) Terphenyl-d14	12.83	244	1332384	78.04	ng	-0.05

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.22	88	160140	40.87	ng	98
3) Pyridine	2.92	79	379057	43.21	ng	100
4) n-Nitrosodimethylamine	2.88	42	168112	43.59	ng	97
6) Aniline	6.40	93	539925	39.20	ng	# 47
8) 2-Chlorophenol	6.52	128	404024	40.99	ng	94
9) Benzaldehyde	6.28	77	232877	41.23	ng	95
10) Phenol	6.40	94	468822	39.15	ng	# 64
11) bis(2-Chloroethyl)ether	6.48	93	415676	43.84	ng	99
12) 1,3-Dichlorobenzene	6.67	146	441529m	42.27	ng	
13) 1,4-Dichlorobenzene	6.75	146	434541	40.38	ng	93
14) 1,2-Dichlorobenzene	6.90	146	383577	38.72	ng	97
15) Benzyl Alcohol	6.89	79	261090	38.44	ng	99
16) 2,2'-oxybis(1-Chloropropan	7.01	45	453982	39.19	ng	57
17) 2-Methylphenol	7.00	107	306046	39.39	ng	# 95
18) Hexachloroethane	7.24	117	163160	41.22	ng	95
19) n-Nitroso-di-n-propylamine	7.16	70	258160	39.68	ng	98
20) 3+4-Methylphenols	7.16	107	376234	39.81	ng	# 74
22) Acetophenone	7.15	105	537156	38.41	ng	# 93
24) Nitrobenzene	7.32	77	393461	35.65	ng	99
25) Isophorone	7.56	82	796944	40.02	ng	99
26) 2-Nitrophenol	7.64	139	238568	45.18	ng	99
27) 2,4-Dimethylphenol	7.69	122	373387	39.37	ng	98
28) bis(2-Chloroethoxy)methane	7.78	93	491334	42.05	ng	99
29) 2,4-Dichlorophenol	7.89	162	306592	38.23	ng	92
30) 1,2,4-Trichlorobenzene	7.96	180	349265	38.80	ng	97
31) Naphthalene	8.04	128	1181183	39.47	ng	99
32) Benzoic acid	7.85	122	226532	32.55	ng	98
33) 4-Chloroaniline	8.10	127	494538	40.91	ng	98
34) Hexachlorobutadiene	8.16	225	194626	40.22	ng	99
35) Caprolactam	8.49	113	104525	41.09	ng	100
36) 4-Chloro-3-methylphenol	8.59	107	342645	38.01	ng	98
37) 2-Methylnaphthalene	8.73	142	659049	33.71	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.90	216	344411	43.01	ng	99
40) Hexachlorocyclopentadiene	8.88	237	73841	34.80	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.01	196	215005	41.81	nq	99
43) 2,4,5-Trichlorophenol	9.06	196	221979	42.52	nq	# 85
45) 1,1'-Biphenyl	9.20	154	892888	38.87	nq	96
46) 2-Chloronaphthalene	9.22	162	650739	39.08	nq	93
47) 2-Nitroaniline	9.32	65	191574	39.32	nq	90
48) Acenaphthylene	9.63	152	968338	36.30	nq	100
49) Dimethylphthalate	9.50	163	816785	41.36	nq	100
50) 2,6-Dinitrotoluene	9.57	165	200630	48.02	nq	# 73
51) Acenaphthene	9.80	154	573529	36.15	nq	98
52) 3-Nitroaniline	9.74	138	224711	44.63	nq	# 78
53) 2,4-Dinitrophenol	9.86	184	66646	34.21	nq	# 73
54) Dibenzofuran	9.97	168	772930	36.89	nq	98
55) 4-Nitrophenol	9.92	139	122314	41.20	nq	90
56) 2,4-Dinitrotoluene	9.97	165	192901	36.54	nq	# 39
57) Fluorene	10.32	166	642331	35.89	nq	100
58) 2,3,4,6-Tetrachlorophenol	10.10	232	156332	40.22	nq	97
59) Diethylphthalate	10.20	149	855512	42.92	nq	98
60) 4-Chlorophenyl-phenylether	10.30	204	331214	39.73	nq	93
61) 4-Nitroaniline	10.36	138	218270	44.02	nq	91
62) Azobenzene	10.46	77	772606	40.47	nq	95
64) 4,6-Dinitro-2-methylphenol	10.38	198	107204	34.35	nq	93
65) n-Nitrosodiphenylamine	10.43	169	574983	35.70	nq	100
66) 4-Bromophenyl-phenylether	10.80	248	202545	38.52	nq	93
67) Hexachlorobenzene	10.86	284	208273	38.31	nq	# 83
68) Atrazine	10.97	200	220753	42.42	nq	100
69) Pentachlorophenol	11.07	266	87224	34.23	nq	98
70) Phenanthrene	11.28	178	1015860	36.16	nq	99
71) Anthracene	11.33	178	1113518	37.83	nq	99
72) Carbazole	11.49	167	968868	38.30	nq	98
73) Di-n-butylphthalate	11.81	149	1113270	35.51	nq	99
74) Fluoranthene	12.46	202	1137681	38.93	nq	93
76) Benzidine	12.59	184	495371	34.99	nq	99
77) Pyrene	12.69	202	1142223	40.39	nq	99
79) Butylbenzylphthalate	13.31	149	544880	42.79	nq	92
80) Benzo(a)anthracene	13.88	228	900831	38.08	nq	99
81) 3,3'-Dichlorobenzidine	13.85	252	637211	36.88	nq	99
82) Chrysene	13.92	228	841160	36.91	nq	99
83) Bis(2-ethylhexyl)phthalate	13.87	149	694776	41.11	nq	# 93
84) Di-n-octyl phthalate	14.48	149	1111916	41.20	nq	100
85) Indeno(1,2,3-cd)pyrene	16.63	276	678618	36.71	nq	99
87) Benzo(b)fluoranthene	14.90	252	969738m	50.08	nq	
88) Benzo(k)fluoranthene	14.92	252	576732m	33.64	nq	
89) Benzo(a)pyrene	15.23	252	706671	41.19	nq	# 97
90) Dibenzo(a,h)anthracene	16.63	278	572041	41.44	nq	98
91) Benzo(g,h,i)perylene	17.03	276	557705	41.75	nq	94

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

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