

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100516\
 Data File : BF090421.D
 Acq On : 6 Oct 2016 11:59
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

SOHIL
 10/6/2016 5:28:02 PM

Quant Time: Oct 06 14:00:35 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF100516.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 05 14:57:00 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.00	152	258019	20.00	ng	0.00
21) Naphthalene-d8	8.29	136	1043555	20.00	ng	0.00
38) Acenaphthene-d10	10.05	164	504901	20.00	ng	0.00
63) Phenanthrene-d10	11.55	188	783392	20.00	ng	0.00
75) Chrysene-d12	14.20	240	534688	20.00	ng	0.00
86) Perylene-d12	15.70	264	465589	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.61	112	1319106	76.30	ng	0.00
7) Phenol-d6	6.62	99	1709954	75.52	ng	0.00
23) Nitrobenzene-d5	7.57	82	1623854	75.69	ng	0.00
41) 2,4,6-Tribromophenol	10.85	330	325464	79.27	ng	0.00
44) 2-Fluorobiphenyl	9.38	172	2550780	84.17	ng	0.00
78) Terphenyl-d14	13.14	244	2088144	87.80	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.69	88	314971	36.96	ng	# 91
3) Pyridine	3.47	79	902613	37.98	ng	# 82
4) n-Nitrosodimethylamine	3.41	42	344736	35.16	ng	# 59
6) Aniline	6.67	93	1160456	37.10	ng	95
8) 2-Chlorophenol	6.78	128	683960	35.95	ng	87
9) Benzaldehyde	6.54	77	463819	40.52	ng	88
10) Phenol	6.65	94	1014411	38.53	ng	91
11) bis(2-Chloroethyl)ether	6.74	93	762002	37.81	ng	95
12) 1,3-Dichlorobenzene	6.94	146	682876	36.01	ng	# 92
13) 1,4-Dichlorobenzene	7.02	146	759120	39.41	ng	97
14) 1,2-Dichlorobenzene	7.18	146	634093m	34.17	ng	
15) Benzyl Alcohol	7.14	79	614407	35.69	ng	96
16) 2,2'-oxybis(1-Chloropropan	7.27	45	1211377	35.44	ng	95
17) 2-Methylphenol	7.25	107	548223	35.04	ng	99
18) Hexachloroethane	7.53	117	266272	35.06	ng	# 89
19) n-Nitroso-di-n-propylamine	7.42	70	625193	37.65	ng	# 81
20) 3+4-Methylphenols	7.41	107	778288	38.43	ng	# 73
22) Acetophenone	7.41	105	875644	35.43	ng	# 98
24) Nitrobenzene	7.58	77	784509	36.83	ng	93
25) Isophorone	7.83	82	1495511	38.13	ng	97
26) 2-Nitrophenol	7.90	139	331867	36.20	ng	# 73
27) 2,4-Dimethylphenol	7.94	122	608877	34.15	ng	95
28) bis(2-Chloroethoxy)methane	8.04	93	953084	41.09	ng	97
29) 2,4-Dichlorophenol	8.14	162	485503	35.82	ng	# 85
30) 1,2,4-Trichlorobenzene	8.23	180	563517	40.34	ng	99
31) Naphthalene	8.31	128	1788718	35.53	ng	99
32) Benzoic acid	8.06	122	514934	41.55	ng	96
33) 4-Chloroaniline	8.36	127	807803	38.82	ng	96
34) Hexachlorobutadiene	8.44	225	314628	40.31	ng	99
35) Caprolactam	8.74	113	170998	37.52	ng	91
36) 4-Chloro-3-methylphenol	8.84	107	653381	38.50	ng	99
37) 2-Methylnaphthalene	9.01	142	1242243	40.76	ng	98
39) 1,2,4,5-Tetrachlorobenzene	9.17	216	476802	35.82	ng	# 98
40) Hexachlorocyclopentadiene	9.16	237	156263	21.42	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.29	196	407294	44.34	nq	98
43) 2,4,5-Trichlorophenol	9.32	196	326139	37.55	nq	93
45) 1,1'-Biphenyl	9.48	154	1487732	39.76	nq	93
46) 2-Chloronaphthalene	9.50	162	1183209	42.80	nq	95
47) 2-Nitroaniline	9.59	65	432011	39.61	nq #	79
48) Acenaphthylene	9.91	152	1659006	36.37	nq	99
49) Dimethylphthalate	9.78	163	1371752	42.42	nq	98
50) 2,6-Dinitrotoluene	9.83	165	314236	43.77	nq #	83
51) Acenaphthene	10.09	154	1015100	35.69	nq	98
52) 3-Nitroaniline	10.01	138	365866	43.63	nq #	84
53) 2,4-Dinitrophenol	10.10	184	86001	26.87	nq #	1
54) Dibenzofuran	10.26	168	1324299	35.98	nq	92
55) 4-Nitrophenol	10.15	139	303339	41.34	nq	88
56) 2,4-Dinitrotoluene	10.23	165	378302	39.60	nq #	40
57) Fluorene	10.60	166	1130890	36.56	nq	99
58) 2,3,4,6-Tetrachlorophenol	10.37	232	233023m	35.92	nq	
59) Diethylphthalate	10.47	149	1319622	39.44	nq	93
60) 4-Chlorophenyl-phenylether	10.60	204	580693	41.28	nq #	86
61) 4-Nitroaniline	10.61	138	311599	36.88	nq	93
62) Azobenzene	10.76	77	1558509	40.33	nq	88
64) 4,6-Dinitro-2-methylphenol	10.65	198	147522	30.17	nq #	54
65) n-Nitrosodiphenylamine	10.71	169	1081581	41.20	nq	97
66) 4-Bromophenyl-phenylether	11.09	248	331910	43.22	nq #	87
67) Hexachlorobenzene	11.16	284	349369	40.85	nq #	88
68) Atrazine	11.24	200	299167	46.22	nq	98
69) Pentachlorophenol	11.34	266	192910	37.90	nq	98
70) Phenanthrene	11.57	178	1613359	40.22	nq	98
71) Anthracene	11.62	178	1463385	35.72	nq	98
72) Carbazole	11.78	167	1503031	39.49	nq	96
73) Di-n-butylphthalate	12.11	149	2020188	43.62	nq #	97
74) Fluoranthene	12.76	202	1343161	34.14	nq	98
76) Benzidine	12.87	184	571492	40.65	nq	98
77) Pyrene	12.99	202	1309244	36.72	nq	99
79) Butylbenzylphthalate	13.61	149	683764	37.38	nq #	69
80) Benzo(a)anthracene	14.18	228	1200913	40.03	nq	98
81) 3,3'-Dichlorobenzidine	14.14	252	450152	41.39	nq #	97
82) Chrysene	14.22	228	1182604	42.83	nq	98
83) Bis(2-ethylhexyl)phthalate	14.18	149	1019913	39.22	nq #	97
84) Di-n-octyl phthalate	14.80	149	1683984	43.68	nq #	92
85) Indeno(1,2,3-cd)pyrene	17.22	276	1080173	54.87	nq #	94
87) Benzo(b)fluoranthene	15.25	252	1224739	41.30	nq #	94
88) Benzo(k)fluoranthene	15.29	252	862556m	30.96	nq	
89) Benzo(a)pyrene	15.63	252	987470	38.63	nq #	94
90) Dibenzo(a,h)anthracene	17.24	278	906022	41.66	nq #	94
91) Benzo(g,h,i)perylene	17.69	276	844679	40.47	nq	99

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

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