

Data Path : Z:\HPCHEM1\BNA F\DATA\BF092616\
 Data File : BF090239.D
 Acq On : 27 Sep 2016 1:12
 Operator : UM/SJ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040

Manual Integrations
 APPROVED

sohil
 9/27/2016 1:06:35 PM

Quant Time: Sep 27 01:49:44 2016
 Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092016.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 26 15:58:38 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.48	152	139239	20.00	ng	0.00
21) Naphthalene-d8	7.77	136	537767	20.00	ng	0.00
38) Acenaphthene-d10	9.52	164	262668	20.00	ng	0.01
63) Phenanthrene-d10	10.98	188	385348	20.00	ng	0.00
75) Chrysene-d12	13.60	240	255713	20.00	ng	0.00
86) Perylene-d12	14.91	264	261833	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.03	112	679431	79.54	ng	0.00
7) Phenol-d6	6.15	99	860605	81.58	ng	0.01
23) Nitrobenzene-d5	7.06	82	761920	82.13	ng	0.00
41) 2,4,6-Tribromophenol	10.30	330	168196	74.64	ng	0.00
44) 2-Fluorobiphenyl	8.86	172	1282813	95.88	ng	0.00
78) Terphenyl-d14	12.56	244	768797	76.33	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	1.84	88	162643	35.09	ng	97
3) Pyridine	2.42	79	419770	41.76	ng	97
4) n-Nitrosodimethylamine	2.39	42	121190	40.40	ng	98
6) Aniline	6.14	93	520212	39.56	ng	# 70
8) 2-Chlorophenol	6.26	128	373315	37.98	ng	93
9) Benzaldehyde	6.02	77	204854	44.56	ng	90
10) Phenol	6.16	94	538120	43.15	ng	# 62
11) bis(2-Chloroethyl)ether	6.23	93	348729	35.86	ng	99
12) 1,3-Dichlorobenzene	6.42	146	403588	39.30	ng	95
13) 1,4-Dichlorobenzene	6.50	146	416825	39.76	ng	97
14) 1,2-Dichlorobenzene	6.65	146	351785m	37.67	ng	
15) Benzyl Alcohol	6.64	79	230816	36.69	ng	99
16) 2,2'-oxybis(1-Chloropropan	6.78	45	451781	38.87	ng	76
17) 2-Methylphenol	6.76	107	291065	37.60	ng	98
18) Hexachloroethane	6.99	117	129224	34.56	ng	94
19) n-Nitroso-di-n-propylamine	6.92	70	263925	42.59	ng	97
20) 3+4-Methylphenols	6.92	107	371369	39.95	ng	95
22) Acetophenone	6.91	105	517846	39.93	ng	96
24) Nitrobenzene	7.07	77	334535	34.61	ng	99
25) Isophorone	7.32	82	736497	38.00	ng	99
26) 2-Nitrophenol	7.39	139	188913	39.69	ng	99
27) 2,4-Dimethylphenol	7.45	122	308162	36.44	ng	97
28) bis(2-Chloroethoxy)methane	7.54	93	426201	36.35	ng	99
29) 2,4-Dichlorophenol	7.64	162	307182	41.41	ng	96
30) 1,2,4-Trichlorobenzene	7.71	180	288211	38.90	ng	100
31) Naphthalene	7.79	128	1021091	39.88	ng	100
32) Benzoic acid	7.60	122	222107	38.26	ng	97
33) 4-Chloroaniline	7.85	127	388427	36.57	ng	99
34) Hexachlorobutadiene	7.92	225	157312	40.25	ng	98
35) Caprolactam	8.24	113	93179	37.95	ng	96
36) 4-Chloro-3-methylphenol	8.35	107	349034	41.64	ng	92
37) 2-Methylnaphthalene	8.48	142	597725	37.54	ng	98
39) 1,2,4,5-Tetrachlorobenzene	8.65	216	260066	43.14	ng	99
40) Hexachlorocyclopentadiene	8.64	237	29186	9.81	ng	98

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	8.76	196	181704	42.29	nq	97
43) 2,4,5-Trichlorophenol	8.81	196	193326	43.46	nq	96
45) 1,1'-Biphenyl	8.95	154	736315	39.18	nq	98
46) 2-Chloronaphthalene	8.97	162	623138	44.95	nq	97
47) 2-Nitroaniline	9.07	65	194746	46.59	nq	98
48) Acenaphthylene	9.37	152	861523	38.26	nq	99
49) Dimethylphthalate	9.27	163	723245	43.23	nq	99
50) 2,6-Dinitrotoluene	9.31	165	145168	38.93	nq	97
51) Acenaphthene	9.55	154	543529	40.67	nq	98
52) 3-Nitroaniline	9.48	138	202507	46.32	nq	89
53) 2,4-Dinitrophenol	9.59	184	18264	11.56	nq	# 76
54) Dibenzofuran	9.71	168	697787	39.68	nq	94
55) 4-Nitrophenol	9.66	139	104613	34.93	nq	# 77
56) 2,4-Dinitrotoluene	9.71	165	161643	36.66	nq	94
57) Fluorene	10.06	166	540142	40.75	nq	99
58) 2,3,4,6-Tetrachlorophenol	9.84	232	122206	36.69	nq	99
59) Diethylphthalate	9.95	149	645445	39.95	nq	99
60) 4-Chlorophenyl-phenylether	10.06	204	230707	38.17	nq	95
61) 4-Nitroaniline	10.09	138	187068	41.98	nq	90
62) Azobenzene	10.22	77	675145	40.14	nq	97
64) 4,6-Dinitro-2-methylphenol	10.12	198	33208	13.76	nq	96
65) n-Nitrosodiphenylamine	10.18	169	564323	44.53	nq	100
66) 4-Bromophenyl-phenylether	10.54	248	154952	40.86	nq	# 72
67) Hexachlorobenzene	10.60	284	184972	41.85	nq	# 86
68) Atrazine	10.72	200	136301	39.25	nq	94
69) Pentachlorophenol	10.80	266	92441	36.85	nq	96
70) Phenanthrene	11.00	178	719187	39.91	nq	99
71) Anthracene	11.05	178	755185	39.95	nq	99
72) Carbazole	11.22	167	763813	38.22	nq	97
73) Di-n-butylphthalate	11.56	149	919614	39.58	nq	99
74) Fluoranthene	12.18	202	681578	35.23	nq	92
76) Benzidine	12.32	184	338072	39.50	nq	99
77) Pyrene	12.40	202	638177	36.47	nq	99
79) Butylbenzylphthalate	13.05	149	360739	42.15	nq	93
80) Benzo(a)anthracene	13.59	228	561484	40.81	nq	99
81) 3,3'-Dichlorobenzidine	13.56	252	246226	43.40	nq	# 95
82) Chrysene	13.62	228	550937	41.11	nq	99
83) Bis(2-ethylhexyl)phthalate	13.61	149	470091	42.92	nq	99
84) Di-n-octyl phthalate	14.22	149	843324	44.69	nq	99
85) Indeno(1,2,3-cd)pyrene	16.06	276	522210	48.19	nq	97
87) Benzo(b)fluoranthene	14.57	252	516411m	35.62	nq	
88) Benzo(k)fluoranthene	14.61	252	574037m	42.19	nq	
89) Benzo(a)pyrene	14.87	252	524505	38.46	nq	99
90) Dibenzo(a,h)anthracene	16.08	278	441269	41.46	nq	99
91) Benzo(g,h,i)perylene	16.40	276	413073	39.65	nq	97

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

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