

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021016\
 Data File : BF085008.D
 Acq On : 10 Feb 2016 22:18
 Operator : SJ/IZ
 Sample : SSTDCCC040
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 SSTDCCC040EC

Manual Integrations
 APPROVED

UMANGI
 2/11/2016 10:47:02 AM

Quant Time: Feb 11 01:05:35 2016
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 10 13:28:51 2016
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.63	152	201904	20.00	ng	0.00
21) Naphthalene-d8	7.92	136	782688	20.00	ng	0.00
38) Acenaphthene-d10	9.67	164	327499	20.00	ng	0.00
63) Phenanthrene-d10	11.14	188	493036	20.00	ng	0.00
75) Chrysene-d12	13.77	240	350666	20.00	ng	0.00
86) Perylene-d12	15.13	264	347397	20.00	ng	0.00

System Monitoring Compounds

5) 2-Fluorophenol	5.21	112	953700	73.57	ng	0.01
7) Phenol-d6	6.30	99	1180938	73.49	ng	0.01
23) Nitrobenzene-d5	7.21	82	1131753	81.46	ng	0.01
41) 2,4,6-Tribromophenol	10.46	330	238846	69.92	ng	0.00
44) 2-Fluorobiphenyl	8.99	172	1743669	92.34	ng	0.00
78) Terphenyl-d14	12.73	244	1261942	81.55	ng	0.00

Target Compounds

						Qvalue
2) 1,4-Dioxane	2.08	88	224717	39.58	ng	# 76
3) Pyridine	2.72	79	626301	42.34	ng	# 70
4) n-Nitrosodimethylamine	2.68	42	293559	38.37	ng	80
6) Aniline	6.28	93	744777	37.88	ng	# 91
8) 2-Chlorophenol	6.42	128	561748	38.29	ng	91
9) Benzaldehyde	6.17	77	336718	35.48	ng	98
10) Phenol	6.31	94	735602	40.86	ng	97
11) bis(2-Chloroethyl)ether	6.38	93	575850	41.02	ng	88
12) 1,3-Dichlorobenzene	6.57	146	599004	38.64	ng	98
13) 1,4-Dichlorobenzene	6.65	146	589767m	38.06	ng	
14) 1,2-Dichlorobenzene	6.80	146	527857	36.57	ng	98
15) Benzyl Alcohol	6.79	79	391404	35.27	ng	90
16) 2,2'-oxybis(1-Chloropropan	6.92	45	875305	37.07	ng	98
17) 2-Methylphenol	6.91	107	407813	35.14	ng	# 87
18) Hexachloroethane	7.14	117	164971	27.64	ng	# 82
19) n-Nitroso-di-n-propylamine	7.06	70	347485	34.50	ng	# 74
20) 3+4-Methylphenols	7.07	107	550354	38.21	ng	91
22) Acetophenone	7.05	105	778286	37.73	ng	# 99
24) Nitrobenzene	7.22	77	515881	34.67	ng	95
25) Isophorone	7.47	82	1022899	37.86	ng	95
26) 2-Nitrophenol	7.54	139	318298	43.38	ng	94
27) 2,4-Dimethylphenol	7.60	122	512554	40.16	ng	90
28) bis(2-Chloroethoxy)methane	7.68	93	562727	35.11	ng	99
29) 2,4-Dichlorophenol	7.79	162	423465	38.77	ng	95
30) 1,2,4-Trichlorobenzene	7.86	180	460454	37.33	ng	96
31) Naphthalene	7.94	128	1510731	38.48	ng	99
32) Benzoic acid	7.75	122	332467	40.72	ng	94
33) 4-Chloroaniline	8.00	127	622659	38.35	ng	97
34) Hexachlorobutadiene	8.06	225	243277	34.20	ng	97
35) Caprolactam	8.40	113	134205	39.87	ng	# 78
36) 4-Chloro-3-methylphenol	8.50	107	465585	37.37	ng	91
37) 2-Methylnaphthalene	8.63	142	815043	33.17	ng	97
39) 1,2,4,5-Tetrachlorobenzene	8.80	216	455491	43.79	ng	98
40) Hexachlorocyclopentadiene	8.78	237	6182	3.85	ng	94

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42) 2,4,6-Trichlorophenol	8.91	196	258686	38.61	ng	96
43) 2,4,5-Trichlorophenol	8.96	196	261373	38.39	ng #	86
45) 1,1'-Biphenyl	9.10	154	1111768	38.01	ng	99
46) 2-Chloronaphthalene	9.12	162	843311	39.15	ng	93
47) 2-Nitroaniline	9.22	65	262613	38.50	ng	96
48) Acenaphthylene	9.53	152	1250054	38.24	ng	98
49) Dimethylphthalate	9.40	163	944017	37.85	ng #	90
50) 2,6-Dinitrotoluene	9.47	165	218520	40.10	ng #	87
51) Acenaphthene	9.70	154	793546	37.31	ng	97
52) 3-Nitroaniline	9.63	138	208726	34.09	ng	93
53) 2,4-Dinitrophenol	9.75	184	38535	20.01	ng	89
54) Dibenzofuran	9.87	168	970773	37.35	ng	96
55) 4-Nitrophenol	9.82	139	141659	31.48	ng #	83
56) 2,4-Dinitrotoluene	9.87	165	265637	38.22	ng	92
57) Fluorene	10.22	166	846137	38.98	ng	98
58) 2,3,4,6-Tetrachlorophenol	10.00	232	199329	38.46	ng	87
59) Diethylphthalate	10.10	149	925718	38.01	ng	99
60) 4-Chlorophenyl-phenylether	10.20	204	359013	38.28	ng	88
61) 4-Nitroaniline	10.25	138	241832	40.53	ng #	82
62) Azobenzene	10.36	77	819499	34.99	ng	88
64) 4,6-Dinitro-2-methylphenol	10.27	198	65759	21.61	ng #	42
65) n-Nitrosodiphenylamine	10.33	169	671840	42.91	ng	98
66) 4-Bromophenyl-phenylether	10.70	248	241411	44.33	ng #	87
67) Hexachlorobenzene	10.75	284	233443	39.34	ng #	76
68) Atrazine	10.87	200	224087	45.33	ng	96
69) Pentachlorophenol	10.96	266	106319	38.13	ng	99
70) Phenanthrene	11.16	178	981021	35.87	ng	97
71) Anthracene	11.22	178	1150628	41.76	ng	100
72) Carbazole	11.38	167	987679	41.11	ng	99
73) Di-n-butylphthalate	11.72	149	1274904	41.68	ng	99
74) Fluoranthene	12.35	202	1024515	37.02	ng	93
76) Benzidine	12.48	184	532615	43.72	ng	98
77) Pyrene	12.57	202	993285	37.00	ng	99
79) Butylbenzylphthalate	13.21	149	485824	39.89	ng #	95
80) Benzo(a)anthracene	13.76	228	871365	40.04	ng	99
81) 3,3'-Dichlorobenzidine	13.74	252	360197	46.74	ng #	99
82) Chrysene	13.81	228	880260	41.71	ng	99
83) Bis(2-ethylhexyl)phthalate	13.77	149	587209	38.74	ng	98
84) Di-n-octyl phthalate	14.39	149	1060799	42.42	ng #	90
85) Indeno(1,2,3-cd)pyrene	16.40	276	758681	45.67	ng	98
87) Benzo(b)fluoranthene	14.77	252	814365m	34.89	ng	
88) Benzo(k)fluoranthene	14.79	252	764798m	39.63	ng	
89) Benzo(a)pyrene	15.09	252	780203	39.23	ng	99
90) Dibenzo(a,h)anthracene	16.40	278	636699	38.03	ng	98
91) Benzo(g,h,i)perylene	16.77	276	619448	36.73	ng	100

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

