

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061715\
 Data File : BF080040.D
 Acq On : 17 Jun 2015 23:54
 Operator : TP/IZ
 Sample : G2651-09MS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NC-HC-001MS

Manual Integrations
 APPROVED

apatel
 6/18/2015 4:07:22 PM

Quant Time: Jun 18 01:29:26 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 17 16:41:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.34	152	43894	20.00	ng	0.00
21) Naphthalene-d8	8.63	136	172406	20.00	ng	0.00
38) Acenaphthene-d10	10.40	164	85857	20.00	ng	0.00
63) Phenanthrene-d10	11.90	188	166175	20.00	ng	-0.01
75) Chrysene-d12	14.79	240	192277	20.00	ng	-0.07
86) Perylene-d12	16.63	264	194116	20.00	ng	-0.11

System Monitoring Compounds

5) 2-Fluorophenol	5.96	112	355377	143.32	ng	0.01
7) Phenol-d6	6.94	99	466725	141.44	ng	0.00
23) Nitrobenzene-d5	7.90	82	285075	96.26	ng	0.00
41) 2,4,6-Tribromophenol	11.19	330	128344	152.87	ng	-0.01
44) 2-Fluorobiphenyl	9.70	172	597221	99.20	ng	0.00
78) Terphenyl-d14	13.56	244	735300	89.31	ng	-0.03

Target Compounds

						Qvalue
2) 1,4-Dioxane	3.24	88	36301	30.80	ng	87
3) Pyridine	4.06	79	89416	28.53	ng	# 83
4) n-Nitrosodimethylamine	3.96	42	57162	38.37	ng	92
6) Aniline	7.00	93	180609	39.78	ng	95
8) 2-Chlorophenol	7.12	128	136935	47.78	ng	92
9) Benzaldehyde	6.88	77	88726	46.58	ng	# 74
10) Phenol	6.96	94	151857	42.17	ng	90
11) bis(2-Chloroethyl)ether	7.05	93	127475	44.62	ng	99
12) 1,3-Dichlorobenzene	7.28	146	150957m	43.85	ng	
13) 1,4-Dichlorobenzene	7.36	146	143467	43.82	ng	98
14) 1,2-Dichlorobenzene	7.51	146	139113m	43.66	ng	
15) Benzyl Alcohol	7.46	79	127163	47.13	ng	# 79
16) 2,2'-oxybis(1-Chloropropan	7.60	45	192386	45.37	ng	89
17) 2-Methylphenol	7.57	107	109975	44.92	ng	# 84
18) Hexachloroethane	7.86	117	46771	42.90	ng	# 72
19) n-Nitroso-di-n-propylamine	7.73	70	96293	42.03	ng	# 80
20) 3+4-Methylphenols	7.72	107	144949	45.88	ng	86
22) Acetophenone	7.74	105	194628	48.44	ng	# 83
24) Nitrobenzene	7.91	77	139071	45.50	ng	# 63
25) Isophorone	8.15	82	273858	45.95	ng	# 87
26) 2-Nitrophenol	8.23	139	59499	46.50	ng	# 32
27) 2,4-Dimethylphenol	8.25	122	118079	44.26	ng	# 76
28) bis(2-Chloroethoxy)methane	8.36	93	170153	48.59	ng	# 93
29) 2,4-Dichlorophenol	8.47	162	107689	47.13	ng	88
30) 1,2,4-Trichlorobenzene	8.57	180	116151	44.77	ng	96
31) Naphthalene	8.65	128	402704m	44.68	ng	
32) Benzoic acid	8.32	122	61829	38.33	ng	# 71
33) 4-Chloroaniline	8.69	127	133210m	34.53	ng	
34) Hexachlorobutadiene	8.77	225	70048	43.84	ng	99
35) Caprolactam	9.03	113	33582	45.28	ng	91
36) 4-Chloro-3-methylphenol	9.16	107	123535	47.94	ng	# 79
37) 2-Methylnaphthalene	9.35	142	275476	47.25	ng	# 93
39) 1,2,4,5-Tetrachlorobenzene	9.51	216	115497	46.23	ng	99
40) Hexachlorocyclopentadiene	9.51	237	46870	39.86	ng	99

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF061715\
 Data File : BF080040.D
 Acq On : 17 Jun 2015 23:54
 Operator : TP/IZ
 Sample : G2651-09MS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_F
 ClientSampleId :
 NC-HC-001MS

Manual Integrations
 APPROVED

apatel
 6/18/2015 4:07:22 PM

Quant Time: Jun 18 01:29:26 2015
 Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 17 16:41:53 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
42) 2,4,6-Trichlorophenol	9.62	196	81689	48.05	ng	99
43) 2,4,5-Trichlorophenol	9.66	196	95420	48.72	ng	97
45) 1,1'-Biphenyl	9.81	154	328486	47.02	ng	94
46) 2-Chloronaphthalene	9.84	162	269036	47.47	ng	96
47) 2-Nitroaniline	9.92	65	79411	51.04	ng	# 68
48) Acenaphthylene	10.26	152	438450	46.35	ng	98
49) Dimethylphthalate	10.09	163	343115	53.16	ng	# 97
50) 2,6-Dinitrotoluene	10.16	165	65929	50.95	ng	# 81
51) Acenaphthene	10.44	154	273177	47.20	ng	96
52) 3-Nitroaniline	10.33	138	69121	46.30	ng	# 73
53) 2,4-Dinitrophenol	10.44	184	39666	75.65	ng	# 1
54) Dibenzofuran	10.61	168	391967	47.73	ng	# 91
55) 4-Nitrophenol	10.48	139	122306	102.49	ng	# 82
56) 2,4-Dinitrotoluene	10.57	165	78017	47.51	ng	# 87
57) Fluorene	10.95	166	295001	45.27	ng	92
58) 2,3,4,6-Tetrachlorophenol	10.72	232	83209	50.32	ng	93
59) Diethylphthalate	10.80	149	321332	49.44	ng	99
60) 4-Chlorophenyl-phenylether	10.94	204	150307	48.00	ng	97
61) 4-Nitroaniline	10.95	138	70343	45.62	ng	# 68
62) Azobenzene	11.10	77	317750	48.03	ng	91
64) 4,6-Dinitro-2-methylphenol	10.98	198	31379	40.07	ng	# 71
65) n-Nitrosodiphenylamine	11.05	169	268634	49.64	ng	93
66) 4-Bromophenyl-phenylether	11.43	248	86826	49.14	ng	93
67) Hexachlorobenzene	11.51	284	90725	46.20	ng	# 81
68) Atrazine	11.57	200	85661	50.11	ng	82
69) Pentachlorophenol	11.70	266	109450	98.36	ng	98
70) Phenanthrene	11.93	178	459530	45.68	ng	97
71) Anthracene	11.98	178	475142	49.05	ng	97
72) Carbazole	12.13	167	430291	46.52	ng	95
73) Di-n-butylphthalate	12.45	149	525934	49.22	ng	# 98
74) Fluoranthene	13.17	202	533416	49.78	ng	89
76) Benzidine	13.28	184	273095	50.81	ng	98
77) Pyrene	13.42	202	549598	46.42	ng	96
79) Butylbenzylphthalate	14.08	149	245777	51.70	ng	88
80) Benzo(a)anthracene	14.78	228	539834	46.09	ng	96
81) 3,3'-Dichlorobenzidine	14.72	252	200171	49.55	ng	# 95
82) Chrysene	14.83	228	510385	47.25	ng	94
83) Bis(2-ethylhexyl)phthalate	14.75	149	371450	49.44	ng	# 92
84) Di-n-octyl phthalate	15.52	149	658065	51.11	ng	95
85) Indeno(1,2,3-cd)pyrene	18.48	276	625952	45.95	ng	# 89
87) Benzo(b)fluoranthene	16.09	252	522442	44.45	ng	# 86
88) Benzo(k)fluoranthene	16.13	252	551067m	46.04	ng	
89) Benzo(a)pyrene	16.55	252	531450	47.61	ng	# 87
90) Dibenzo(a,h)anthracene	18.51	278	523361	45.81	ng	# 83
91) Benzo(g,h,i)perylene	19.05	276	512019	44.39	ng	# 75

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

