

Data Path : S:\HPCHEM1\BNA_L\data\BL072612\
 Data File : BL000657.D
 Acq On : 26 Jul 2012 11:15
 Operator : HLM
 Sample : SSTD050
 Misc : SSTD050 HLM
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_L
 LabSampleId :
 SSTD050

Quant Time: Sep 06 06:42:24 2012
 Quant Method : H:\VOL1\GCMSDATA\L\METHODS\BL072512.M
 Quant Title : SW-846 Method 8270
 QLast Update : Thu Jul 26 11:02:15 2012
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
1 I	1,4-Dichlorobenzene-d4	40.000	40.000	0.0	89	0.01
2 T	1,4-Dioxane	50.000	0.000	100.0#	0	-4.08#
3 T	N-Nitrosodimethylamine	50.000	55.000	-10.0	95	0.00
4 T	Pyridine	50.000	56.213	-12.4	95	0.00
5 S	2-Fluorophenol	50.000	53.588	-7.2	91	0.01
6 T	Benzaldehyde	50.000	110.915	-121.8#	100	0.01
7 T	Aniline	50.000	52.604	-5.2	88	0.00
8 T	bis(2-Chloroethyl)ether	50.000	52.383	-4.8	90	0.01
9 S	Phenol-d5	50.000	52.903	-5.8	88	0.01
10 MC	Phenol	50.000	53.632	-7.3	88	0.01
11 M	2-Chlorophenol	50.000	51.197	-2.4	85	0.01
12 T	1,3-Dichlorobenzene	50.000	51.534	-3.1	88	0.01
13 MC	1,4-Dichlorobenzene	50.000	51.777	-3.6	89	0.00
14 T	1,2-Dichlorobenzene	50.000	52.606	-5.2	88	0.01
15 T	Benzyl alcohol	50.000	53.315	-6.6	86	0.01
16 T	2,2-oxybis(1-chloropropane)	50.000	55.828	-11.7	95	0.01
17 T	2-Methylphenol	50.000	52.288	-4.6	88	0.01
18 T	Acetophenone	50.000	50.308	-0.6	87	0.01
19 T	Hexachloroethane	50.000	50.763	-1.5	86	0.00
20 MP	N-Nitroso-di-n-propylamine	50.000	52.543	-5.1	90	0.01
21 T	4-Methylphenol	50.000	52.572	-5.1	88	0.01
22 I	Naphthalene-d8	40.000	40.000	0.0	87	0.01
23 S	Nitrobenzene-d5	50.000	52.724	-5.4	89	0.01
24 T	Nitrobenzene	50.000	53.581	-7.2	89	0.00
25 T	Isophorone	50.000	51.535	-3.1	86	0.01
26 TC	2-Nitrophenol	50.000	51.825	-3.7	83	0.00
27 T	2,4-Dimethylphenol	50.000	52.579	-5.2	87	0.01
28 T	Benzoic Acid	100.000	91.568	8.4	75	-0.02
29 T	bis(2-Chloroethoxy)methane	50.000	52.175	-4.3	86	0.01
30 TC	2,4-Dichlorophenol	50.000	50.903	-1.8	81	0.01
31 T	2,6-Dichlorophenol	50.000	0.000	100.0#	0	-12.90#
32 M	1,2,4-Trichlorobenzene	50.000	50.872	-1.7	83	0.01
33 T	Naphthalene	50.000	52.875	-5.8	87	0.01
34 T	4-Chloroaniline	50.000	52.607	-5.2	84	0.01
35 TC	Hexachlorobutadiene	50.000	52.522	-5.0	85	0.01
36 T	Caprolactam	50.000	50.057	-0.1	79	0.00
37 MC	4-Chloro-3-methylphenol	50.000	52.095	-4.2	83	0.01
38 T	2-Methylnaphthalene	50.000	52.659	-5.3	86	0.01
39 I	Acenaphthene-d10	40.000	40.000	0.0	82	0.01
40 TP	Hexachlorocyclopentadiene	50.000	48.350	3.3	78	0.01
41 TC	2,4,6-Trichlorophenol	50.000	51.199	-2.4	81	0.01
42 T	2,4,5-Trichlorophenol	50.000	50.153	-0.3	79	0.01
43 S	2-Fluorobiphenyl	50.000	51.587	-3.2	83	0.01
44 T	1,1-Biphenyl	50.000	51.713	-3.4	84	0.00

Data Path : S:\HPCHEM1\BNA_L\data\BL072612\
 Data File : BL000657.D
 Acq On : 26 Jul 2012 11:15
 Operator : HLM
 Sample : SSTD050
 Misc : SSTD050 HLM
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_L
 LabSampleId :
 SSTD050

Quant Time: Sep 06 06:42:24 2012
 Quant Method : H:\VOL1\GCMSDATA\L\METHODS\BL072512.M
 Quant Title : SW-846 Method 8270
 QLast Update : Thu Jul 26 11:02:15 2012
 Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
 Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
45 T	2-Chloronaphthalene	50.000	51.355	-2.7	83	0.00
46 T	2-Nitroaniline	50.000	53.036	-6.1	83	0.01
47 T	Acenaphthylene	50.000	51.395	-2.8	81	0.01
48 T	Dimethylphthalate	50.000	49.219	1.6	78	0.01
49 T	2,6-Dinitrotoluene	50.000	49.256	1.5	77	0.00
50 MC	Acenaphthene	50.000	52.255	-4.5	84	0.01
51 T	3-Nitroaniline	50.000	53.170	-6.3	77	0.01
52 TP	2,4-Dinitrophenol	50.000	43.126	13.7	67	0.01
53 T	Dibenzofuran	50.000	50.845	-1.7	80	0.01
54 M	2,4-Dinitrotoluene	50.000	50.458	-0.9	78	0.01
55 MP	4-Nitrophenol	50.000	47.469	5.1	74	0.01
56 T	2,3,4,6-Tetrachlorophenol	50.000	50.153	-0.3	77	0.01
57 T	Fluorene	50.000	52.322	-4.6	82	0.00
58 T	4-Chlorophenyl-phenylether	50.000	51.428	-2.9	81	0.01
59 T	Diethylphthalate	50.000	48.293	3.4	75	0.00
60 T	4-Nitroaniline	50.000	49.865	0.3	77	0.00
61 S	2,4,6-Tribromophenol	50.000	51.420	-2.8	79	0.00
62 I	Phenanthrene-d10	40.000	40.000	0.0	77	0.01
63 T	1,2,4,5-Tetrachlorobenzene	50.000	55.915	-11.8	85	0.01
64 T	4,6-Dinitro-2-methylphenol	50.000	46.936	6.1	70	0.00
65 TC	n-Nitrosodiphenylamine	50.000	55.039	-10.1	79	0.01
66 T	1,2-Diphenylhydrazine	50.000	55.629	-11.3	82	0.01
67 T	4-Bromophenyl-phenylether	50.000	54.036	-8.1	77	0.01
68 T	Hexachlorobenzene	50.000	54.536	-9.1	80	0.01
69 T	Atrazine	50.000	48.492	3.0	73	0.01
70 MC	Pentachlorophenol	50.000	45.306	9.4	67	0.01
71 T	Pentachloronitrophenol	50.000	0.000	100.0#	0	-18.46#
72 T	Benzidine	50.000	0.000	100.0#	0	0.01
73 T	Phenanthrene	50.000	52.053	-4.1	78	0.01
74 T	Anthracene	50.000	52.869	-5.7	77	0.00
75 T	Carbazole	50.000	50.151	-0.3	76	0.00
76 T	Di-n-butylphthalate	50.000	47.508	5.0	73	0.00
77 TC	Fluoranthene	50.000	50.087	-0.2	81	0.01
78 I	Chrysene-d12	40.000	40.000	0.0	102	0.00
79 M	Pyrene	50.000	45.373	9.3	83	0.01
80 S	Terphenyl-d14	50.000	43.387	13.2	82	0.01
81 T	Butylbenzylphthalate	50.000	47.372	5.3	92	0.01
82 T	3,3-Dichlorobenzidine	50.000	57.943	-15.9	106	0.00
83 T	Benzo[a]anthracene	50.000	50.589	-1.2	98	0.01
84 T	Chrysene	50.000	53.044	-6.1	103	0.00
85 T	bis(2-Ethylhexyl)phthalate	50.000	48.880	2.2	92	0.00
86 I	Perylene-d12	40.000	40.000	0.0	118	0.01
87 TC	Di-n-octylphthalate	50.000	46.569	6.9	101	0.00

Data Path : S:\HPCHEM1\BNA_L\data\BL072612\
Data File : BL000657.D
Acq On : 26 Jul 2012 11:15
Operator : HLM
Sample : SSTD050
Misc : SSTD050 HLM
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_L
LabSampleId :
SSTD050

Quant Time: Sep 06 06:42:24 2012
Quant Method : H:\VOL1\GCMSDATA\L\METHODS\BL072512.M
Quant Title : SW-846 Method 8270
QLast Update : Thu Jul 26 11:02:15 2012
Response via : Initial Calibration

Min. RRF : 0.050 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	Amount	Calc.	%Dev	Area%	Dev(min)
88 T	Benzo[b]fluoranthene	50.000	48.476	3.0	116	0.01
89 T	Benzo[k]fluoranthene	50.000	54.232	-8.5	117	0.01
90 TC	Benzo[a]pyrene	50.000	51.173	-2.3	115	0.01
91 T	Indeno[1,2,3-cd]pyrene	50.000	53.633	-7.3	116	0.02
92 T	Dibenz[a,h]anthracene	50.000	52.830	-5.7	116	0.02
93 T	Benzo[g,h,i]perylene	50.000	52.735	-5.5	117	0.02

(#) = Out of Range

SPCC's out = 0 CCC's out = 0