Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_L\data\BL072612\
Data File : BL000659.D
Acq On : 26 Jul 2012 12:29
Operator : HLM
Sample : SSTD050ICV
Misc : SSTD050ICV HLM
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_L
ClientSampleId : ICVBL072512

Quant Time: Sep 06 06:45:36 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	102	0.01
2	T	1,4-Dioxane	0.000	0.000#	0.0	0#	-4.08#
3	T	N-Nitrosodimethylamine	1.286	1.245	3.2	96	0.09
4	T	Pyridine	1.963	1.872	4.6	93	0.08
5	S	2-Fluorophenol	1.254	0.000#	100.0#	0#	-7.51#
6	T	Benzaldehyde	0.289	1.138	-293.8#	205#	0.02
7	T	Aniline	1.872	1.284	31.4#	66	0.00
8	T	bis(2-Chloroethyl)ether	1.255	1.197	4.6	95	0.01
9	S	Phenol-d5	1.448	0.000#	100.0#	0#	-9.34#
10	MC	Phenol	1.675	1.620	3.3	91	0.01
11	M	2-Chlorophenol	1.035	0.988	4.5	92	0.01
12	T	1,3-Dichlorobenzene	1.048	1.000	4.6	94	0.01
13	MC	1,4-Dichlorobenzene	1.047	1.002	4.3	94	0.01
14	T	1,2-Dichlorobenzene	0.958	0.928	3.1	94	0.01
15	T	Benzyl alcohol	0.671	0.648	3.4	90	0.01
16	T	2,2-oxybis(1-chloropropane)	1.943	1.888	2.8	95	0.01
17	T	2-Methylphenol	0.963	0.898	6.7	90	0.02
18	T	Acetophenone	1.128	2.031	-80.1#	180#	0.01
19	T	Hexachloroethane	0.406	0.378	6.9	91	0.00
20	MP	N-Nitroso-di-n-propylamine	0.724	0.648	10.5	88	0.00
21	T	4-Methylphenol	0.902	0.824	8.6	88	0.01
22	I	Naphthalene-d8	1.000	1.000	0.0	100	0.01
23	S	Nitrobenzene-d5	0.304	0.024#	92.1#	8#	-0.07
24	T	Nitrobenzene	0.316	0.293	7.3	89	0.00
25	T	Isophorone	0.608	0.574	5.6	91	0.01
26	TC	2-Nitrophenol	0.160	0.153	4.4	89	0.00
27	T	2,4-Dimethylphenol	0.242	0.233	3.7	92	0.01
28	T	Benzoic Acid	0.169	0.170	-0.6	86	0.00
29	T	bis(2-Chloroethoxy)methane	0.362	0.350	3.3	92	0.01
30	TC	2,4-Dichlorophenol	0.195	0.189	3.1	90	0.01
31	T	2,6-Dichlorophenol	0.000	0.000#	0.0	0#	-12.90#
32	M	1,2,4-Trichlorobenzene	0.215	0.207	3.7	91	0.01
33	T	Naphthalene	0.690	0.663	3.9	91	0.00
34	T	4-Chloroaniline	0.292	0.155	46.9#	48#	0.00
35	TC	Hexachlorobutadiene	0.092	0.089	3.3	90	0.00
36	T	Caprolactam	0.078	0.000#	100.0#	0#	-13.13#
37	MC	4-Chloro-3-methylphenol	0.182	0.182	0.0	92	0.00
38	T	2-Methylnaphthalene	0.384	0.373	2.9	92	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.01
40	TP	Hexachlorocyclopentadiene	0.223	0.179	19.7	79	0.01
41	TC	2,4,6-Trichlorophenol	0.282	0.264	6.4	90	0.01
42	T	2,4,5-Trichlorophenol	0.308	0.293	4.9	91	0.00
43	S	2-Fluorobiphenyl	0.959	0.000#	100.0#	0#	-14.30#
44	T	1,1-Biphenyl	1.086	1.835	-69.0#	168#	0.00

Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_L\data\BL072612\
Data File : BL000659.D
Acq On : 26 Jul 2012 12:29
Operator : HLM
Sample : SSTD050ICV
Misc : SSTD050ICV HLM
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_L
ClientSampleId : ICVBL072512

Quant Time: Sep 06 06:45:36 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	T	2-Chloronaphthalene	0.887	0.800	9.8	89	0.00
46	T	2-Nitroaniline	0.272	0.260	4.4	91	0.01
47	T	Acenaphthylene	1.279	1.170	8.5	88	0.00
48	T	Dimethylphthalate	0.877	0.815	7.1	90	0.00
49	T	2,6-Dinitrotoluene	0.231	0.213	7.8	87	0.00
50	MC	Acenaphthene	0.759	0.724	4.6	93	0.00
51	T	3-Nitroaniline	0.222	0.151	32.0#	60	0.00
52	TP	2,4-Dinitrophenol	0.111	0.100	9.9	77	0.00
53	T	Dibenzofuran	1.008	0.963	4.5	92	0.01
54	M	2,4-Dinitrotoluene	0.256	0.249	2.7	92	0.00
55	MP	4-Nitrophenol	0.169	0.161	4.7	86	0.00
56	T	2,3,4,6-Tetrachlorophenol	0.187	0.178	4.8	90	0.00
57	T	Fluorene	0.720	0.700	2.8	93	0.00
58	T	4-Chlorophenyl-phenylether	0.321	0.313	2.5	92	0.01
59	T	Diethylphthalate	0.728	0.691	5.1	89	0.00
60	T	4-Nitroaniline	0.161	0.140	13.0	77	0.00
61	S	2,4,6-Tribromophenol	0.085	0.000#	100.0#	0#	-17.03#
62	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.01
63	T	1,2,4,5-Tetrachlorobenzene	0.338	0.586	-73.4#	175#	0.01
64	T	4,6-Dinitro-2-methylphenol	0.110	0.106	3.6	89	0.00
65	TC	n-Nitrosodiphenylamine	0.507	0.483	4.7	91	0.00
66	T	1,2-Diphenylhydrazine	0.766	0.741	3.3	94	0.01
67	T	4-Bromophenyl-phenylether	0.147	0.141	4.1	91	0.01
68	T	Hexachlorobenzene	0.155	0.151	2.6	94	0.01
69	T	Atrazine	0.111	0.000#	100.0#	0#	-17.82#
70	MC	Pentachlorophenol	0.076	0.071	6.6	83	0.01
71	T	Pentachloronitrophenol	0.000	0.000#	0.0	0#	-18.46#
72	T	Benzidine	0.000	0.000#	0.0	0#	-20.33#
73	T	Phenanthrene	0.677	0.641	5.3	94	0.00
74	T	Anthracene	0.697	0.665	4.6	92	0.00
75	T	Carbazole	0.598	0.564	5.7	94	0.00
76	T	Di-n-butylphthalate	0.793	0.724	8.7	93	0.00
77	TC	Fluoranthene	0.573	0.524	8.6	99	0.00
78	I	Chrysene-d12	1.000	1.000	0.0	112	0.00
79	M	Pyrene	0.956	0.939	1.8	99	0.00
80	S	Terphenyl-d14	0.521	0.000#	100.0#	0#	-20.65#
81	T	Butylbenzylphthalate	0.453	0.418	7.7	98	0.01
82	T	3,3-Dichlorobenzidine	0.162	0.000#	100.0#	0#	-21.85#
83	T	Benzo[a]anthracene	0.724	0.683	5.7	100	0.01
84	T	Chrysene	0.686	0.679	1.0	105	0.00
85	T	bis(2-Ethylhexyl)phthalate	0.662	0.588	11.2	92	0.00
86	I	Perylene-d12	1.000	1.000	0.0	112	0.01
87	TC	Di-n-octylphthalate	1.347	1.177	12.6	91	0.00

Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_L\data\BL072612\
Data File : BL000659.D
Acq On : 26 Jul 2012 12:29
Operator : HLM
Sample : SSTD050ICV
Misc : SSTD050ICV HLM
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_L
ClientSampleId : ICVBL072512

Quant Time: Sep 06 06:45:36 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	T	Benzo[b]fluoranthene	0.844	0.782	7.3	105	0.01
89	T	Benzo[k]fluoranthene	0.779	0.734	5.8	96	0.01
90	TC	Benzo[a]pyrene	0.775	0.706	8.9	98	0.01
91	T	Indeno[1,2,3-cd]pyrene	0.763	0.728	4.6	98	0.02
92	T	Dibenz[a,h]anthracene	0.638	0.594	6.9	97	0.01
93	T	Benzo[g,h,i]perylene	0.665	0.632	5.0	100	0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 0