Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021915\
Data File : BG016015.D
Acq On : 19 Feb 2015 22:51
Operator : TP/IZ
Sample : SSTDCCC020
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Feb 20 06:19:12 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-MA-BG021915.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Feb 20 05:38:39 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% 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	95	0.00
2		Benzaldehyde	1.203	1.190	1.1	95	0.00
3	S	Phenol-d5	1.903	1.853	2.6	94	0.00
4		Phenol	2.006	1.973	1.6	96	0.00
5	S	Bis-(2-Chloroethyl)ether-d8	1.124	1.090	3.0	96	0.00
6		Bis(2-Chloroethyl)ether	1.577	1.558	1.2	97	0.00
7	S	2-Chlorophenol-d4	1.408	1.399	0.6	97	0.00
8		2-Chlorophenol	1.414	1.408	0.4	97	0.00
9		2-Methylphenol	1.427	1.377	3.5	94	0.00
10		2,2'-oxybis(1-Chloropropane	2.005	1.921	4.2	96	0.00
11	S	4-Methylphenol-d8	1.408	1.389	1.3	96	0.00
12		Acetophenone	2.133	2.085	2.3	97	0.00
13	P	N-Nitroso-di-n-propylamine	1.202	1.162	3.3	95	0.00
14		4-Methylphenol	1.578	1.560	1.1	96	0.00
15		Hexachloroethane	0.561	0.544	3.0	95	0.00
16	I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
17	S	Nitrobenzene-d5	0.141	0.141	0.0	100	0.00
18		Nitrobenzene	0.331	0.339	-2.4	103	0.00
19		Isophorone	0.720	0.687	4.6	95	0.00
20	S	2-Nitrophenol-d4	0.130	0.129	0.8	101	0.00
21	C	2-Nitrophenol	0.147	0.149	-1.4	104	0.00
22		2,4-Dimethylphenol	0.353	0.342	3.1	96	0.00
23		Bis(2-Chloroethoxy)methane	0.449	0.434	3.3	96	0.00
24	S	2,4-Dichlorophenol-d3	0.283	0.276	2.5	97	0.00
25	C	2,4-Dichlorophenol	0.277	0.273	1.4	97	0.00
26		Naphthalene	1.068	1.038	2.8	96	0.00
27	S	4-Chloroaniline-d4	0.312	0.314	-0.6	103	0.00
28		4-Chloroaniline	0.304	0.306	-0.7	102	0.00
29	C	Hexachlorobutadiene	0.155	0.147	5.2	96	0.00
30		Caprolactam	0.129	0.124	3.9	93	0.00
31	C	4-Chloro-3-methylphenol	0.316	0.306	3.2	95	0.00
32		2-Methylnaphthalene	0.750	0.730	2.7	97	0.00
33	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.00
34		1,2,4,5-Tetrachlorobenzene	0.537	0.512	4.7	96	0.00
35		Hexachlorocyclopentadiene	0.238	0.236	0.8	98	0.00
36	C	2,4,6-Trichlorophenol	0.308	0.298	3.2	97	0.00
37		2,4,5-Trichlorophenol	0.344	0.338	1.7	97	0.00
38		1,1'-Biphenyl	1.547	1.522	1.6	98	0.00
39		2-Chloronaphthalene	1.163	1.133	2.6	97	0.00
40		2-Nitroaniline	0.296	0.299	-1.0	104	0.00
41	S	Dimethylphthalate-d6	1.337	1.327	0.7	96	0.00
42		Dimethylphthalate	1.394	1.379	1.1	97	0.00
43		2,6-Dinitrotoluene	0.241	0.241	0.0	101	0.00
44	S	Acenaphthylene-d8	1.892	1.853	2.1	96	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021915\
Data File : BG016015.D
Acq On : 19 Feb 2015 22:51
Operator : TP/IZ
Sample : SSTDCCC020
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Feb 20 06:19:12 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-MA-BG021915.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Feb 20 05:38:39 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		Acenaphthylene	2.006	1.983	1.1	97	0.00
46		3-Nitroaniline	0.270	0.278	-3.0	102	0.00
47	C	Acenaphthene	1.331	1.306	1.9	97	0.00
48		2,4-Dinitrophenol	0.088	0.068	22.7#	92	0.00
49	S	4-Nitrophenol-d4	0.269	0.266	1.1	103	0.00
50		4-Nitrophenol	0.155	0.155	0.0	100	0.00
51		Dibenzofuran	1.783	1.751	1.8	98	0.00
52		2,4-Dinitrotoluene	0.345	0.355	-2.9	100	0.00
53		2,3,4,6-Tetrachlorophenol	0.277	0.281	-1.4	101	0.00
54		Diethylphthalate	1.412	1.426	-1.0	99	0.00
55	S	Fluorene-d10	1.344	1.309	2.6	97	0.00
56		Fluorene	1.555	1.528	1.7	98	0.00
57		4-Chlorophenyl-phenylether	0.702	0.671	4.4	96	0.00
58		4-Nitroaniline	0.368	0.387	-5.2	103	0.00
59	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
60	S	4,6-Dinitro-2-methylphenol-	0.077	0.070	9.1	103	0.00
61		4,6-Dinitro-2-methylphenol	0.083	0.075	9.6	100	0.00
62		N-Nitrosodiphenylamine	0.625	0.610	2.4	98	0.00
63		4-Bromophenyl-phenylether	0.205	0.194	5.4	96	0.00
64		Hexachlorobenzene	0.232	0.224	3.4	98	0.00
65		Atrazine	0.201	0.201	0.0	99	0.00
66	C	Pentachlorophenol	0.098	0.090	8.2	103	0.00
67		Phenanthrene	1.112	1.100	1.1	98	0.00
68	S	Anthracene-d10	0.949	0.931	1.9	98	0.00
69		Anthracene	1.135	1.134	0.1	98	0.00
70		1,2,3,4-Tetrachlorobenzene	0.254	0.239	5.9	96	0.00
71		Pentachlorobenzene	0.249	0.234	6.0	94	0.00
72		Carbazole	1.071	1.076	-0.5	99	0.00
73		Di-n-butylphthalate	1.145	1.223	-6.8	101	0.00
74	C	Fluoranthene	1.217	1.260	-3.5	102	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	101	0.00
76	S	Pyrene-d10	0.939	0.923	1.7	101	0.00
77		Pyrene	1.217	1.215	0.2	101	0.00
78		Butylbenzylphthalate	0.476	0.495	-4.0	103	0.00
79		3,3'-Dichlorobenzidine	0.274	0.257	6.2	94	0.00
80		Benzo(a)anthracene	1.112	1.111	0.1	100	0.00
81		Bis(2-ethylhexyl)phthalate	0.625	0.653	-4.5	103	0.00
82		Chrysene	1.060	1.062	-0.2	101	0.00
83	I	Perylene-d12	1.000	1.000	0.0	101	0.00
84		Di-n-octyl phthalate	1.058	1.112	-5.1	104	0.00
85		Benzo(b)fluoranthene	1.115	1.084	2.8	99	0.00
86		Benzo(k)fluoranthene	1.120	1.109	1.0	100	0.00
87	S	Benzo(a)pyrene-d12	0.907	0.872	3.9	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG021915\
Data File : BG016015.D
Acq On : 19 Feb 2015 22:51
Operator : TP/IZ
Sample : SSTDCCC020
Misc :
ALS Vial : 4 Sample Multiplier: 1

Quant Time: Feb 20 06:19:12 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM01.2-EPA-MA-BG021915.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Feb 20 05:38:39 2015
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 20% Max. R.T. Dev 0.50min
Max. RRF Dev : 20% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88 C	Benzo(a)pyrene	1.108	1.087	1.9	100	0.00
89	Indeno(1,2,3-cd)pyrene	1.289	1.231	4.5	99	0.00
90	Dibenzo(a,h)anthracene	1.077	1.044	3.1	100	0.00
91	Benzo(g,h,i)perylene	1.077	1.033	4.1	100	0.00

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