Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051816\
Data File : BG021949.D
Acq On : 19 May 2016 1:57
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTD02001

Quant Time: May 19 03:58:00 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG051816.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 19 03:03:18 2016
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	129	0.00
2	S	1,4-Dioxane-d8	0.436	0.411	5.7	117	0.00
3	S	Phenol-d5	1.766	1.814	-2.7	135	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	0.941	0.957	-1.7	128	0.00
5	S	2-Chlorophenol-d4	1.450	1.549	-6.8	131	0.00
6	S	4-Methylphenol-d8	1.573	1.587	-0.9	126	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	129	0.00
8	S	Nitrobenzene-d5	0.140	0.145	-3.6	128	0.00
9	S	2-Nitrophenol-d4	0.150	0.146	2.7	122	0.00
10	S	2,4-Dichlorophenol-d3	0.265	0.271	-2.3	128	0.00
11		Naphthalene	0.973	0.996	-2.4	129	0.00
12	S	4-Chloroaniline-d4	0.208	0.207	0.5	138	0.00
13		2-Methylnaphthalene	0.705	0.706	-0.1	127	0.00
14		1-Methylnaphthalene	0.721	0.718	0.4	124	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	125	0.00
16	S	Dimethylphthalate-d6	1.542	1.466	4.9	118	0.00
17	S	Acenaphthylene-d8	1.986	1.987	-0.1	121	0.00
18		Acenaphthylene	2.068	2.082	-0.7	123	0.00
19	C	Acenaphthene	1.420	1.431	-0.8	124	0.00
20	S	4-Nitrophenol-d4	0.271	0.270	0.4	129	0.00
21	S	Fluorene-d10	1.388	1.351	2.7	121	0.00
22		Fluorene	1.573	1.523	3.2	118	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	120	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.099	0.097	2.0	126	0.00
25		Phenanthrene	1.069	1.086	-1.6	120	0.00
26	S	Anthracene-d10	0.921	0.918	0.3	119	0.00
27		Anthracene	1.074	1.089	-1.4	118	0.00
28	C	Fluoranthene	1.111	1.172	-5.5	124	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	129	0.00
30	S	Pyrene-d10	1.055	1.038	1.6	125	0.00
31		Pyrene	1.377	1.316	4.4	121	0.00
32		Benzo(a)anthracene	1.109	1.118	-0.8	128	0.00
33		Chrysene	1.079	1.130	-4.7	135	0.00
34	I	Perylene-d12	1.000	1.000	0.0	128	0.00
35		Benzo(b)fluoranthene	1.150	1.168	-1.6	129	0.00
36		Benzo(k)fluoranthene	1.131	1.154	-2.0	130	0.00
37	S	Benzo(a)pyrene-d12	0.888	0.896	-0.9	128	0.00
38	C	Benzo(a)pyrene	1.099	1.099	0.0	127	0.00
39		Indeno(1,2,3-cd)pyrene	1.250	1.265	-1.2	130	0.00
40		Dibenzo(a,h)anthracene	1.035	1.027	0.8	125	0.00
41		Benzo(g,h,i)perylene	1.057	1.061	-0.4	127	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG051816\
Data File : BG021949.D
Acq On : 19 May 2016 1:57
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTD02001

Quant Time: May 19 03:58:00 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG051816.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 19 03:03:18 2016
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)

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