Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG051916\
Data File : BG021972.D
Acq On : 19 May 2016 21:14
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTD02005

Quant Time: May 20 08:27:44 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG051816.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 20 08:11:42 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	171#	0.00
2	S	1,4-Dioxane-d8	0.436	0.382	12.4	144	0.00
3	S	Phenol-d5	1.766	1.712	3.1	168#	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	0.941	0.896	4.8	158#	0.00
5	S	2-Chlorophenol-d4	1.450	1.460	-0.7	164#	0.00
6	S	4-Methylphenol-d8	1.573	1.497	4.8	157#	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	160#	0.00
8	S	Nitrobenzene-d5	0.140	0.146	-4.3	159#	0.00
9	S	2-Nitrophenol-d4	0.150	0.154	-2.7	160#	0.00
10	S	2,4-Dichlorophenol-d3	0.265	0.274	-3.4	161#	0.00
11		Naphthalene	0.973	0.968	0.5	156#	0.00
12	S	4-Chloroaniline-d4	0.208	0.215	-3.4	178#	0.00
13		2-Methylnaphthalene	0.705	0.695	1.4	155#	0.00
14		1-Methylnaphthalene	0.721	0.683	5.3	146	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	145	0.00
16	S	Dimethylphthalate-d6	1.542	1.424	7.7	133	0.00
17	S	Acenaphthylene-d8	1.986	1.999	-0.7	142	0.00
18		Acenaphthylene	2.068	2.073	-0.2	142	0.00
19	C	Acenaphthene	1.420	1.409	0.8	141	0.00
20	S	4-Nitrophenol-d4	0.271	0.226	16.6	126	0.00
21	S	Fluorene-d10	1.388	1.344	3.2	140	0.00
22		Fluorene	1.573	1.506	4.3	136	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	132	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.099	0.086	13.1	124	0.00
25		Phenanthrene	1.069	1.077	-0.7	131	0.00
26	S	Anthracene-d10	0.921	0.912	1.0	130	0.00
27		Anthracene	1.074	1.083	-0.8	130	0.00
28	C	Fluoranthene	1.111	1.116	-0.5	130	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	134	0.00
30	S	Pyrene-d10	1.055	1.046	0.9	131	0.00
31		Pyrene	1.377	1.310	4.9	125	0.00
32		Benzo(a)anthracene	1.109	1.121	-1.1	133	0.00
33		Chrysene	1.079	1.046	3.1	130	0.00
34	I	Perylene-d12	1.000	1.000	0.0	134	0.00
35		Benzo(b)fluoranthene	1.150	1.152	-0.2	133	0.00
36		Benzo(k)fluoranthene	1.131	1.129	0.2	134	0.00
37	S	Benzo(a)pyrene-d12	0.888	0.900	-1.4	136	0.00
38	C	Benzo(a)pyrene	1.099	1.092	0.6	132	0.00
39		Indeno(1,2,3-cd)pyrene	1.250	1.229	1.7	133	0.00
40		Dibenzo(a,h)anthracene	1.035	1.065	-2.9	137	0.00
41		Benzo(g,h,i)perylene	1.057	1.028	2.7	129	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG051916\
Data File : BG021972.D
Acq On : 19 May 2016 21:14
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTD02005

Quant Time: May 20 08:27:44 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG051816.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 20 08:11:42 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