Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM062616\
Data File : BM006063.D
Acq On : 27 Jun 2016 05:22
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02037

Quant Time: Jun 27 07:05:18 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 27 05:15:31 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	215#	0.00
2	S	1,4-Dioxane-d8	0.448	0.421	6.0	197#	0.00
3	S	Phenol-d5	2.104	1.894	10.0	196#	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	1.170	1.101	5.9	198#	0.00
5	S	2-Chlorophenol-d4	1.465	1.418	3.2	207#	0.00
6	S	4-Methylphenol-d8	1.776	1.519	14.5	185#	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	192#	0.00
8	S	Nitrobenzene-d5	0.152	0.155	-2.0	193#	0.00
9	S	2-Nitrophenol-d4	0.169	0.171	-1.2	189#	0.00
10	S	2,4-Dichlorophenol-d3	0.325	0.327	-0.6	190#	0.00
11		Naphthalene	0.999	1.015	-1.6	192#	0.00
12	S	4-Chloroaniline-d4	0.353	0.397	-12.5	172#	0.00
13		2-Methylnaphthalene	0.804	0.770	4.2	178#	0.00
14		1-Methylnaphthalene	0.770	0.728	5.5	176#	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	158#	0.00
16	S	Dimethylphthalate-d6	1.712	1.566	8.5	144	0.00
17	S	Acenaphthylene-d8	1.986	2.046	-3.0	159#	0.00
18		Acenaphthylene	2.094	2.156	-3.0	158#	0.00
19	C	Acenaphthene	1.330	1.362	-2.4	160#	0.00
20	S	4-Nitrophenol-d4	0.365	0.324	11.2	131	0.00
21	S	Fluorene-d10	1.462	1.410	3.6	150	0.00
22		Fluorene	1.602	1.527	4.7	146	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	139	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.111	0.071	36.0#	86	0.00
25		Phenanthrene	1.076	1.110	-3.2	141	0.00
26	S	Anthracene-d10	0.919	0.947	-3.0	139	0.00
27		Anthracene	1.111	1.143	-2.9	139	0.00
28	C	Fluoranthene	1.378	1.326	3.8	128	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
30	S	Pyrene-d10	0.876	1.138	-29.9#	126	0.00
31		Pyrene	1.145	1.479	-29.2#	124	0.00
32		Benzo(a)anthracene	1.188	1.243	-4.6	101	0.00
33		Chrysene	1.085	1.128	-4.0	100	0.00
34	I	Perylene-d12	1.000	1.000	0.0	76	0.00
35		Benzo(b)fluoranthene	1.162	1.249	-7.5	81	0.00
36		Benzo(k)fluoranthene	1.081	1.166	-7.9	79	0.00
37	S	Benzo(a)pyrene-d12	0.903	0.931	-3.1	77	0.00
38	C	Benzo(a)pyrene	1.114	1.151	-3.3	77	0.00
39		Indeno(1,2,3-cd)pyrene	1.328	1.283	3.4	71	-0.02
40		Dibenzo(a,h)anthracene	1.104	1.068	3.3	71	-0.02
41		Benzo(g,h,i)perylene	1.149	1.111	3.3	71	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM062616\
Data File : BM006063.D
Acq On : 27 Jun 2016 05:22
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02037

Quant Time: Jun 27 07:05:18 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 27 05:15:31 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