Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062816\
Data File : BM006092.D
Acq On : 28 Jun 2016 12:31
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jun 28 13:24:25 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 28 02:11:41 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	103	0.00
2	S	1,4-Dioxane-d8	0.431	0.333	22.7#	100	0.00
3	S	Phenol-d5	1.634	1.401	14.3	107	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	0.963	0.841	12.7	105	0.00
5	S	2-Chlorophenol-d4	1.254	1.038	17.2	105	0.00
6	S	4-Methylphenol-d8	1.315	1.132	13.9	107	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
8	S	Nitrobenzene-d5	0.135	0.110	18.5	108	0.00
9	S	2-Nitrophenol-d4	0.154	0.127	17.5	108	0.00
10	S	2,4-Dichlorophenol-d3	0.281	0.234	16.7	109	0.00
11		Naphthalene	0.876	0.719	17.9	108	0.00
12	S	4-Chloroaniline-d4	0.222	0.204	8.1	107	0.00
13		2-Methylnaphthalene	0.665	0.560	15.8	110	0.00
14		1-Methylnaphthalene	0.632	0.528	16.5	110	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
16	S	Dimethylphthalate-d6	1.434	1.156	19.4	111	0.00
17	S	Acenaphthylene-d8	1.740	1.414	18.7	111	0.00
18		Acenaphthylene	1.822	1.475	19.0	109	0.00
19	C	Acenaphthene	1.162	0.937	19.4	108	0.00
20	S	4-Nitrophenol-d4	0.272	0.227	16.5	108	0.00
21	S	Fluorene-d10	1.223	0.990	19.1	111	0.00
22		Fluorene	1.323	1.098	17.0	111	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.071	0.039	45.1#	92	0.00
25		Phenanthrene	0.946	0.766	19.0	110	0.00
26	S	Anthracene-d10	0.802	0.651	18.8	109	0.00
27		Anthracene	0.973	0.797	18.1	110	0.00
28	C	Fluoranthene	1.159	0.959	17.3	110	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
30	S	Pyrene-d10	0.805	0.657	18.4	108	0.00
31		Pyrene	1.055	0.864	18.1	108	0.00
32		Benzo(a)anthracene	1.040	0.847	18.6	105	0.00
33		Chrysene	0.957	0.791	17.3	105	0.00
34	I	Perylene-d12	1.000	1.000	0.0	101	0.00
35		Benzo(b)fluoranthene	1.038	0.821	20.9#	101	0.00
36		Benzo(k)fluoranthene	0.974	0.801	17.8	102	0.00
37	S	Benzo(a)pyrene-d12	0.796	0.637	20.0	101	0.00
38	C	Benzo(a)pyrene	0.981	0.799	18.6	101	0.00
39		Indeno(1,2,3-cd)pyrene	1.019	0.854	16.2	99	0.00
40		Dibenzo(a,h)anthracene	0.843	0.714	15.3	100	0.00
41		Benzo(g,h,i)perylene	0.853	0.706	17.2	99	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062816\
Data File : BM006092.D
Acq On : 28 Jun 2016 12:31
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Quant Time: Jun 28 13:24:25 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 28 02:11:41 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