Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM063016\
Data File : BM006152.D
Acq On : 30 Jun 2016 08:15
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02054

Quant Time: Jun 30 14:47:41 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM062816.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 29 00:57:09 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	146	0.00
2	S	1,4-Dioxane-d8	0.431	0.322	25.3#	137	0.00
3	S	Phenol-d5	1.634	1.331	18.5	144	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	0.963	0.813	15.6	144	0.00
5	S	2-Chlorophenol-d4	1.254	1.006	19.8	144	0.00
6	S	4-Methylphenol-d8	1.315	1.061	19.3	141	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	142	0.00
8	S	Nitrobenzene-d5	0.135	0.111	17.8	146	0.00
9	S	2-Nitrophenol-d4	0.154	0.121	21.4#	138	0.00
10	S	2,4-Dichlorophenol-d3	0.281	0.227	19.2	142	0.00
11		Naphthalene	0.876	0.717	18.2	144	0.00
12	S	4-Chloroaniline-d4	0.222	0.176	20.7#	123	0.00
13		2-Methylnaphthalene	0.665	0.538	19.1	141	0.00
14		1-Methylnaphthalene	0.632	0.508	19.6	142	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	135	0.00
16	S	Dimethylphthalate-d6	1.434	1.120	21.9#	131	0.00
17	S	Acenaphthylene-d8	1.740	1.445	17.0	138	0.00
18		Acenaphthylene	1.822	1.509	17.2	135	0.00
19	C	Acenaphthene	1.162	0.957	17.6	135	0.00
20	S	4-Nitrophenol-d4	0.272	0.216	20.6#	125	0.00
21	S	Fluorene-d10	1.223	0.983	19.6	134	0.00
22		Fluorene	1.323	1.080	18.4	133	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	131	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.071	0.076	-7.0	211#	0.00
25		Phenanthrene	0.946	0.772	18.4	131	0.00
26	S	Anthracene-d10	0.802	0.656	18.2	130	0.00
27		Anthracene	0.973	0.805	17.3	131	0.00
28	C	Fluoranthene	1.159	0.915	21.1	123	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	101	0.00
30	S	Pyrene-d10	0.805	0.768	4.6	122	0.00
31		Pyrene	1.055	1.005	4.7	122	0.00
32		Benzo(a)anthracene	1.040	0.848	18.5	102	0.00
33		Chrysene	0.957	0.778	18.7	100	0.00
34	I	Perylene-d12	1.000	1.000	0.0	75	0.00
35		Benzo(b)fluoranthene	1.038	0.858	17.3	79	0.00
36		Benzo(k)fluoranthene	0.974	0.835	14.3	79	0.00
37	S	Benzo(a)pyrene-d12	0.796	0.641	19.5	75	0.00
38	C	Benzo(a)pyrene	0.981	0.799	18.6	75	0.00
39		Indeno(1,2,3-cd)pyrene	1.019	0.896	12.1	77	-0.01
40		Dibenzo(a,h)anthracene	0.843	0.753	10.7	78	0.00
41		Benzo(g,h,i)perylene	0.853	0.778	8.8	81	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM063016\
Data File : BM006152.D
Acq On : 30 Jun 2016 08:15
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02054

Quant Time: Jun 30 14:47:41 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM062816.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jun 29 00:57:09 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