Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081016\
Data File : BM007001.D
Acq On : 10 Aug 2016 15:03
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02042

Quant Time: Aug 10 18:31:15 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080916.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 10 01:30:54 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	74	0.00
2	S	1,4-Dioxane-d8	0.439	0.433	1.4	71	0.00
3	S	Phenol-d5	1.780	1.718	3.5	70	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	0.995	0.995	0.0	71	0.00
5	S	2-Chlorophenol-d4	1.350	1.334	1.2	72	0.00
6	S	4-Methylphenol-d8	1.507	1.446	4.0	70	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	72	0.00
8	S	Nitrobenzene-d5	0.153	0.149	2.6	69	0.00
9	S	2-Nitrophenol-d4	0.175	0.157	10.3	63	0.00
10	S	2,4-Dichlorophenol-d3	0.342	0.343	-0.3	71	0.00
11		Naphthalene	1.003	1.025	-2.2	72	0.00
12	S	4-Chloroaniline-d4	0.399	0.432	-8.3	71	0.00
13		2-Methylnaphthalene	0.794	0.815	-2.6	72	0.00
14		1-Methylnaphthalene	0.762	0.771	-1.2	71	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	72	0.00
16	S	Dimethylphthalate-d6	1.709	1.715	-0.4	70	0.00
17	S	Acenaphthylene-d8	2.002	2.060	-2.9	71	0.00
18		Acenaphthylene	2.069	2.121	-2.5	71	0.00
19	C	Acenaphthene	1.338	1.359	-1.6	70	0.00
20	S	4-Nitrophenol-d4	0.307	0.292	4.9	67	0.00
21	S	Fluorene-d10	1.453	1.489	-2.5	71	0.00
22		Fluorene	1.588	1.629	-2.6	71	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	71	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.108	0.091	15.7	62	0.00
25		Phenanthrene	1.072	1.117	-4.2	71	0.00
26	S	Anthracene-d10	0.932	0.971	-4.2	71	0.00
27		Anthracene	1.107	1.152	-4.1	71	0.00
28	C	Fluoranthene	1.350	1.413	-4.7	72	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	74	0.00
30	S	Pyrene-d10	0.909	0.909	0.0	72	0.00
31		Pyrene	1.160	1.175	-1.3	72	0.00
32		Benzo(a)anthracene	1.178	1.209	-2.6	75	0.00
33		Chrysene	1.069	1.112	-4.0	75	0.00
34	I	Perylene-d12	1.000	1.000	0.0	86	0.00
35		Benzo(b)fluoranthene	1.242	1.228	1.1	82	0.00
36		Benzo(k)fluoranthene	1.174	1.200	-2.2	84	0.00
37	S	Benzo(a)pyrene-d12	0.919	0.956	-4.0	87	0.00
38	C	Benzo(a)pyrene	1.143	1.178	-3.1	86	0.00
39		Indeno(1,2,3-cd)pyrene	1.133	1.313	-15.9	98	0.00
40		Dibenzo(a,h)anthracene	0.938	1.101	-17.4	99	0.00
41		Benzo(g,h,i)perylene	0.932	1.086	-16.5	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM081016\
Data File : BM007001.D
Acq On : 10 Aug 2016 15:03
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02042

Quant Time: Aug 10 18:31:15 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080916.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 10 01:30:54 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