Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM070116\
Data File : BM006174.D
Acq On : 01 Jul 2016 09:48
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02057

Quant Time: Jul 01 17:31:22 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	140	0.00
2	S	1,4-Dioxane-d8	0.431	0.352	18.3	144	0.00
3	S	Phenol-d5	1.634	1.367	16.3	142	0.01
4	S	Bis-(2-Chloroethyl)ether-d8	0.963	0.833	13.5	141	0.00
5	S	2-Chlorophenol-d4	1.254	1.020	18.7	140	0.00
6	S	4-Methylphenol-d8	1.315	1.115	15.2	143	0.01
7	I	Naphthalene-d8	1.000	1.000	0.0	140	0.00
8	S	Nitrobenzene-d5	0.135	0.111	17.8	144	0.01
9	S	2-Nitrophenol-d4	0.154	0.129	16.2	146	0.01
10	S	2,4-Dichlorophenol-d3	0.281	0.231	17.8	142	0.01
11		Naphthalene	0.876	0.719	17.9	142	0.00
12	S	4-Chloroaniline-d4	0.222	0.172	22.5#	119	0.00
13		2-Methylnaphthalene	0.665	0.551	17.1	143	0.00
14		1-Methylnaphthalene	0.632	0.520	17.7	143	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	142	0.00
16	S	Dimethylphthalate-d6	1.434	1.175	18.1	145	0.00
17	S	Acenaphthylene-d8	1.740	1.418	18.5	142	0.01
18		Acenaphthylene	1.822	1.500	17.7	142	0.01
19	C	Acenaphthene	1.162	0.953	18.0	141	0.00
20	S	4-Nitrophenol-d4	0.272	0.228	16.2	139	0.02
21	S	Fluorene-d10	1.223	0.995	18.6	143	0.00
22		Fluorene	1.323	1.101	16.8	142	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	137	0.01
24	S	4,6-Dinitro-2-methylphenol-	0.071	0.089	-25.4#	258#	0.02
25		Phenanthrene	0.946	0.782	17.3	139	0.01
26	S	Anthracene-d10	0.802	0.663	17.3	137	0.01
27		Anthracene	0.973	0.807	17.1	138	0.01
28	C	Fluoranthene	1.159	0.914	21.1	129	0.01
29	I	Chrysene-d12	1.000	1.000	0.0	104	0.01
30	S	Pyrene-d10	0.805	0.772	4.1	127	0.01
31		Pyrene	1.055	1.012	4.1	126	0.01
32		Benzo(a)anthracene	1.040	0.846	18.7	104	0.01
33		Chrysene	0.957	0.785	18.0	104	0.01
34	I	Perylene-d12	1.000	1.000	0.0	82	0.02
35		Benzo(b)fluoranthene	1.038	0.840	19.1	84	0.02
36		Benzo(k)fluoranthene	0.974	0.837	14.1	87	0.01
37	S	Benzo(a)pyrene-d12	0.796	0.647	18.7	83	0.02
38	C	Benzo(a)pyrene	0.981	0.797	18.8	82	0.02
39		Indeno(1,2,3-cd)pyrene	1.019	0.850	16.6	80	0.03
40		Dibenzo(a,h)anthracene	0.843	0.711	15.7	81	0.03
41		Benzo(g,h,i)perylene	0.853	0.725	15.0	82	0.04

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 0