Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080616\
Data File : BM006973.D
Acq On : 07 Aug 2016 12:53
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02024

Quant Time: Aug 08 03:56:20 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080616.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 03:51:46 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	128	0.00
2	S	1,4-Dioxane-d8	0.438	0.384	12.3	107	0.00
3	S	Phenol-d5	1.896	1.888	0.4	133	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	1.189	1.204	-1.3	128	0.00
5	S	2-Chlorophenol-d4	1.354	1.417	-4.7	132	0.00
6	S	4-Methylphenol-d8	1.583	1.555	1.8	131	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	133	0.00
8	S	Nitrobenzene-d5	0.149	0.157	-5.4	137	0.00
9	S	2-Nitrophenol-d4	0.178	0.184	-3.4	129	0.00
10	S	2,4-Dichlorophenol-d3	0.337	0.343	-1.8	132	0.00
11		Naphthalene	1.027	1.046	-1.9	133	0.00
12	S	4-Chloroaniline-d4	0.403	0.429	-6.5	131	0.00
13		2-Methylnaphthalene	0.822	0.820	0.2	130	0.00
14		1-Methylnaphthalene	0.790	0.778	1.5	128	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	126	0.00
16	S	Dimethylphthalate-d6	1.760	1.712	2.7	120	0.00
17	S	Acenaphthylene-d8	2.036	2.069	-1.6	125	0.00
18		Acenaphthylene	2.136	2.172	-1.7	126	0.00
19	C	Acenaphthene	1.393	1.402	-0.6	126	0.00
20	S	4-Nitrophenol-d4	0.255	0.202	20.8#	120	-0.01
21	S	Fluorene-d10	1.496	1.476	1.3	124	0.00
22		Fluorene	1.761	1.721	2.3	122	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	118	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.083	0.115	-38.6#	227#	0.00
25		Phenanthrene	1.119	1.118	0.1	118	0.00
26	S	Anthracene-d10	0.966	0.976	-1.0	119	0.00
27		Anthracene	1.171	1.170	0.1	118	0.00
28	C	Fluoranthene	1.361	1.377	-1.2	116	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	105	0.00
30	S	Pyrene-d10	0.953	0.981	-2.9	114	0.00
31		Pyrene	1.243	1.272	-2.3	114	0.00
32		Benzo(a)anthracene	1.187	1.209	-1.9	108	0.00
33		Chrysene	1.045	1.057	-1.1	105	0.00
34	I	Perylene-d12	1.000	1.000	0.0	89	0.00
35		Benzo(b)fluoranthene	1.305	1.306	-0.1	93	0.00
36		Benzo(k)fluoranthene	1.260	1.273	-1.0	90	0.00
37	S	Benzo(a)pyrene-d12	0.953	0.953	0.0	88	0.00
38	C	Benzo(a)pyrene	1.184	1.169	1.3	88	0.00
39		Indeno(1,2,3-cd)pyrene	1.201	1.281	-6.7	95	-0.01
40		Dibenzo(a,h)anthracene	1.011	1.079	-6.7	96	-0.01
41		Benzo(g,h,i)perylene	0.946	1.024	-8.2	98	-0.02

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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