Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM070116\
Data File : BM006183.D
Acq On : 01 Jul 2016 16:43
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02058

Quant Time: Jul 01 17:33:05 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	110	0.00
2	S	1,4-Dioxane-d8	0.431	0.352	18.3	113	0.00
3	S	Phenol-d5	1.634	1.650	-1.0	135	0.02
4	S	Bis-(2-Chloroethyl)ether-d8	0.963	0.981	-1.9	131	0.00
5	S	2-Chlorophenol-d4	1.254	1.262	-0.6	136	0.01
6	S	4-Methylphenol-d8	1.315	1.344	-2.2	135	0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	109	0.01
8	S	Nitrobenzene-d5	0.135	0.137	-1.5	139	0.01
9	S	2-Nitrophenol-d4	0.154	0.154	0.0	135	0.01
10	S	2,4-Dichlorophenol-d3	0.281	0.285	-1.4	137	0.02
11		Naphthalene	0.876	0.887	-1.3	137	0.00
12	S	4-Chloroaniline-d4	0.222	0.233	-5.0	125	0.01
13		2-Methylnaphthalene	0.665	0.668	-0.5	135	0.01
14		1-Methylnaphthalene	0.632	0.630	0.3	135	0.01
15	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.01
16	S	Dimethylphthalate-d6	1.434	1.370	4.5	122	0.01
17	S	Acenaphthylene-d8	1.740	1.762	-1.3	128	0.01
18		Acenaphthylene	1.822	1.869	-2.6	128	0.01
19	C	Acenaphthene	1.162	1.177	-1.3	126	0.01
20	S	4-Nitrophenol-d4	0.272	0.252	7.4	111	0.04
21	S	Fluorene-d10	1.223	1.218	0.4	126	0.00
22		Fluorene	1.323	1.343	-1.5	126	0.01
23	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.02
24	S	4,6-Dinitro-2-methylphenol-	0.071	0.072	-1.4	145	0.02
25		Phenanthrene	0.946	0.959	-1.4	118	0.02
26	S	Anthracene-d10	0.802	0.823	-2.6	118	0.02
27		Anthracene	0.973	0.998	-2.6	118	0.02
28	C	Fluoranthene	1.159	1.018	12.2	100	0.02
29	I	Chrysene-d12	1.000	1.000	0.0	59	0.01
30	S	Pyrene-d10	0.805	1.025	-27.3#	95	0.02
31		Pyrene	1.055	1.334	-26.4#	94	0.02
32		Benzo(a)anthracene	1.040	1.049	-0.9	73	0.02
33		Chrysene	0.957	0.958	-0.1	71	0.02
34	I	Perylene-d12	1.000	1.000	0.0	52	0.03
35		Benzo(b)fluoranthene	1.038	1.016	2.1	65	0.02
36		Benzo(k)fluoranthene	0.974	0.991	-1.7	65	0.02
37	S	Benzo(a)pyrene-d12	0.796	0.797	-0.1	65	0.03
38	C	Benzo(a)pyrene	0.981	0.991	-1.0	65	0.03
39		Indeno(1,2,3-cd)pyrene	1.019	1.190	-16.8	72	0.05
40		Dibenzo(a,h)anthracene	0.843	0.992	-17.7	72	0.05
41		Benzo(g,h,i)perylene	0.853	1.027	-20.4#	74	0.06

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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