Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061116\
Data File : BM005726.D
Acq On : 11 Jun 2016 12:40
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02052

Quant Time: Jun 11 23:23:24 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jun 11 05:37:26 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	102	0.00
2	S	1,4-Dioxane-d8	0.461	0.451	2.2	103	0.00
3	S	Phenol-d5	1.726	1.744	-1.0	103	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	1.037	1.079	-4.1	104	0.00
5	S	2-Chlorophenol-d4	1.395	1.443	-3.4	104	0.00
6	S	4-Methylphenol-d8	1.458	1.530	-4.9	106	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
8	S	Nitrobenzene-d5	0.144	0.144	0.0	103	0.00
9	S	2-Nitrophenol-d4	0.163	0.164	-0.6	101	0.00
10	S	2,4-Dichlorophenol-d3	0.296	0.297	-0.3	105	0.00
11		Naphthalene	1.002	0.997	0.5	106	-0.01
12	S	4-Chloroaniline-d4	0.357	0.435	-21.8#	108	0.00
13		2-Methylnaphthalene	0.754	0.747	0.9	104	0.00
14		1-Methylnaphthalene	0.745	0.744	0.1	105	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	106	-0.01
16	S	Dimethylphthalate-d6	1.650	1.651	-0.1	105	0.00
17	S	Acenaphthylene-d8	2.018	2.047	-1.4	106	0.00
18		Acenaphthylene	2.066	2.069	-0.1	105	0.00
19	C	Acenaphthene	1.344	1.353	-0.7	106	0.00
20	S	4-Nitrophenol-d4	0.241	0.241	0.0	104	-0.02
21	S	Fluorene-d10	1.397	1.403	-0.4	105	-0.01
22		Fluorene	1.534	1.562	-1.8	105	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	105	-0.01
24	S	4,6-Dinitro-2-methylphenol-	0.103	0.103	0.0	107	-0.02
25		Phenanthrene	1.095	1.090	0.5	104	-0.01
26	S	Anthracene-d10	0.942	0.965	-2.4	107	0.00
27		Anthracene	1.111	1.133	-2.0	104	0.00
28	C	Fluoranthene	1.286	1.306	-1.6	105	-0.01
29	I	Chrysene-d12	1.000	1.000	0.0	107	-0.01
30	S	Pyrene-d10	0.961	0.910	5.3	105	0.00
31		Pyrene	1.217	1.164	4.4	104	0.00
32		Benzo(a)anthracene	1.155	1.134	1.8	106	0.00
33		Chrysene	1.139	1.120	1.7	105	0.00
34	I	Perylene-d12	1.000	1.000	0.0	111	-0.01
35		Benzo(b)fluoranthene	1.188	1.158	2.5	110	-0.01
36		Benzo(k)fluoranthene	1.201	1.185	1.3	108	-0.01
37	S	Benzo(a)pyrene-d12	0.912	0.911	0.1	111	-0.01
38	C	Benzo(a)pyrene	1.128	1.120	0.7	110	-0.02
39		Indeno(1,2,3-cd)pyrene	1.101	1.148	-4.3	116	-0.02
40		Dibenzo(a,h)anthracene	0.921	0.953	-3.5	115	-0.02
41		Benzo(g,h,i)perylene	0.921	0.938	-1.8	116	-0.03

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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