Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM062716\
Data File : BM006090.D
Acq On : 28 Jun 2016 03:22
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02041

Quant Time: Jun 28 04:18:49 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 28 02:11:41 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	208#	0.00
2	S	1,4-Dioxane-d8	0.448	0.387	13.6	176#	0.00
3	S	Phenol-d5	2.104	1.919	8.8	192#	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	1.170	1.112	5.0	194#	0.00
5	S	2-Chlorophenol-d4	1.465	1.439	1.8	203#	0.00
6	S	4-Methylphenol-d8	1.776	1.580	11.0	187#	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	194#	0.00
8	S	Nitrobenzene-d5	0.152	0.156	-2.6	197#	0.00
9	S	2-Nitrophenol-d4	0.169	0.167	1.2	187#	0.00
10	S	2,4-Dichlorophenol-d3	0.325	0.324	0.3	190#	0.01
11		Naphthalene	0.999	1.006	-0.7	192#	0.00
12	S	4-Chloroaniline-d4	0.353	0.402	-13.9	177#	0.00
13		2-Methylnaphthalene	0.804	0.768	4.5	180#	0.00
14		1-Methylnaphthalene	0.770	0.731	5.1	179#	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	164#	0.00
16	S	Dimethylphthalate-d6	1.712	1.552	9.3	149	0.00
17	S	Acenaphthylene-d8	1.986	2.041	-2.8	166#	0.00
18		Acenaphthylene	2.094	2.138	-2.1	164#	0.00
19	C	Acenaphthene	1.330	1.348	-1.4	165#	0.00
20	S	4-Nitrophenol-d4	0.365	0.319	12.6	134	0.00
21	S	Fluorene-d10	1.462	1.415	3.2	156#	0.00
22		Fluorene	1.602	1.503	6.2	150	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	147	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.111	0.057	48.6#	72	0.00
25		Phenanthrene	1.076	1.101	-2.3	148	0.00
26	S	Anthracene-d10	0.919	0.948	-3.2	148	0.00
27		Anthracene	1.111	1.133	-2.0	147	0.00
28	C	Fluoranthene	1.378	1.330	3.5	136	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
30	S	Pyrene-d10	0.876	1.173	-33.9#	135	0.01
31		Pyrene	1.145	1.500	-31.0#	132	0.01
32		Benzo(a)anthracene	1.188	1.220	-2.7	104	0.00
33		Chrysene	1.085	1.136	-4.7	105	0.00
34	I	Perylene-d12	1.000	1.000	0.0	82	0.00
35		Benzo(b)fluoranthene	1.162	1.261	-8.5	88	0.00
36		Benzo(k)fluoranthene	1.081	1.121	-3.7	82	0.00
37	S	Benzo(a)pyrene-d12	0.903	0.925	-2.4	82	0.00
38	C	Benzo(a)pyrene	1.114	1.134	-1.8	82	0.00
39		Indeno(1,2,3-cd)pyrene	1.328	1.275	4.0	76	0.00
40		Dibenzo(a,h)anthracene	1.104	1.058	4.2	76	0.00
41		Benzo(g,h,i)perylene	1.149	1.109	3.5	76	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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