Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061616\
Data File : BM005822.D
Acq On : 16 Jun 2016 02:14
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02043

Quant Time: Jun 16 04:28:13 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061516.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu Jun 16 01:45:55 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	95	0.00
2	S	1,4-Dioxane-d8	0.432	0.424	1.9	94	0.00
3	S	Phenol-d5	1.643	1.697	-3.3	97	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	0.978	1.010	-3.3	95	0.00
5	S	2-Chlorophenol-d4	1.369	1.407	-2.8	97	0.00
6	S	4-Methylphenol-d8	1.388	1.526	-9.9	101	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	101	0.00
8	S	Nitrobenzene-d5	0.142	0.145	-2.1	100	0.00
9	S	2-Nitrophenol-d4	0.157	0.167	-6.4	103	0.00
10	S	2,4-Dichlorophenol-d3	0.301	0.316	-5.0	105	0.00
11		Naphthalene	0.997	0.982	1.5	102	0.00
12	S	4-Chloroaniline-d4	0.331	0.432	-30.5#	118	0.00
13		2-Methylnaphthalene	0.717	0.761	-6.1	109	0.00
14		1-Methylnaphthalene	0.724	0.763	-5.4	111	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	126	0.00
16	S	Dimethylphthalate-d6	1.588	1.671	-5.2	130	0.00
17	S	Acenaphthylene-d8	2.017	2.035	-0.9	127	0.00
18		Acenaphthylene	2.034	2.030	0.2	127	0.00
19	C	Acenaphthene	1.332	1.338	-0.5	128	0.00
20	S	4-Nitrophenol-d4	0.183	0.226	-23.5#	185#	0.00
21	S	Fluorene-d10	1.322	1.429	-8.1	142	0.00
22		Fluorene	1.431	1.542	-7.8	141	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	166#	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.089	0.110	-23.6#	224#	0.00
25		Phenanthrene	1.097	1.085	1.1	166#	0.00
26	S	Anthracene-d10	0.928	0.926	0.2	167#	0.00
27		Anthracene	1.086	1.097	-1.0	167#	0.00
28	C	Fluoranthene	1.168	1.268	-8.6	177#	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	149	0.00
30	S	Pyrene-d10	0.973	1.091	-12.1	176#	0.00
31		Pyrene	1.228	1.356	-10.4	172#	0.00
32		Benzo(a)anthracene	1.157	1.168	-1.0	150#	0.00
33		Chrysene	1.134	1.115	1.7	149	0.00
34	I	Perylene-d12	1.000	1.000	0.0	115	0.00
35		Benzo(b)fluoranthene	1.195	1.279	-7.0	124	0.00
36		Benzo(k)fluoranthene	1.106	1.174	-6.1	125	0.00
37	S	Benzo(a)pyrene-d12	0.924	0.941	-1.8	119	0.00
38	C	Benzo(a)pyrene	1.141	1.146	-0.4	118	0.00
39		Indeno(1,2,3-cd)pyrene	1.309	1.228	6.2	109	0.00
40		Dibenzo(a,h)anthracene	1.095	1.022	6.7	107	0.00
41		Benzo(g,h,i)perylene	1.134	1.040	8.3	106	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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