Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG052216\
Data File : BG022025.D
Acq On : 22 May 2016 23:24
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTD02011

Quant Time: May 24 11:18:45 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG051816.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri May 20 08:11:42 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	170#	0.00
2	S	1,4-Dioxane-d8	0.436	0.434	0.5	163#	0.00
3	S	Phenol-d5	1.766	1.846	-4.5	181#	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	0.941	0.967	-2.8	170#	0.00
5	S	2-Chlorophenol-d4	1.450	1.563	-7.8	175#	0.00
6	S	4-Methylphenol-d8	1.573	1.571	0.1	164#	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	156#	0.00
8	S	Nitrobenzene-d5	0.140	0.157	-12.1	167#	0.00
9	S	2-Nitrophenol-d4	0.150	0.165	-10.0	167#	0.00
10	S	2,4-Dichlorophenol-d3	0.265	0.303	-14.3	173#	0.00
11		Naphthalene	0.973	1.052	-8.1	165#	0.00
12	S	4-Chloroaniline-d4	0.208	0.369	-77.4#	298#	0.00
13		2-Methylnaphthalene	0.705	0.750	-6.4	163#	0.00
14		1-Methylnaphthalene	0.721	0.756	-4.9	158#	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	143	0.00
16	S	Dimethylphthalate-d6	1.542	1.514	1.8	139	0.00
17	S	Acenaphthylene-d8	1.986	2.138	-7.7	149	0.00
18		Acenaphthylene	2.068	2.169	-4.9	147	0.00
19	C	Acenaphthene	1.420	1.485	-4.6	147	0.00
20	S	4-Nitrophenol-d4	0.271	0.257	5.2	141	0.00
21	S	Fluorene-d10	1.388	1.439	-3.7	147	0.00
22		Fluorene	1.573	1.571	0.1	140	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	125	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.099	0.105	-6.1	142	0.00
25		Phenanthrene	1.069	1.163	-8.8	133	0.00
26	S	Anthracene-d10	0.921	0.990	-7.5	133	0.00
27		Anthracene	1.074	1.148	-6.9	130	0.00
28	C	Fluoranthene	1.111	1.236	-11.3	135	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	131	0.00
30	S	Pyrene-d10	1.055	1.110	-5.2	135	0.00
31		Pyrene	1.377	1.402	-1.8	130	0.00
32		Benzo(a)anthracene	1.109	1.194	-7.7	138	0.00
33		Chrysene	1.079	1.087	-0.7	131	0.00
34	I	Perylene-d12	1.000	1.000	0.0	129	0.00
35		Benzo(b)fluoranthene	1.150	1.245	-8.3	138	0.00
36		Benzo(k)fluoranthene	1.131	1.218	-7.7	138	0.00
37	S	Benzo(a)pyrene-d12	0.888	0.997	-12.3	144	0.00
38	C	Benzo(a)pyrene	1.099	1.200	-9.2	139	0.00
39		Indeno(1,2,3-cd)pyrene	1.250	1.274	-1.9	131	0.00
40		Dibenzo(a,h)anthracene	1.035	1.130	-9.2	139	0.00
41		Benzo(g,h,i)perylene	1.057	0.984	6.9	118	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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