Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG052516\
Data File : BG022070.D
Acq On : 25 May 2016 23:25
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTD02017

Quant Time: May 26 03:45:32 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\SOM02.2-EPA-BG051816.M
Quant Title : SVOA CALIBRATION
QLast Update : Thu May 26 03:11:46 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	132	0.00
2	S	1,4-Dioxane-d8	0.436	0.330	24.3#	96	0.00
3	S	Phenol-d5	1.766	1.805	-2.2	137	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	0.941	0.926	1.6	126	0.00
5	S	2-Chlorophenol-d4	1.450	1.493	-3.0	130	0.00
6	S	4-Methylphenol-d8	1.573	1.545	1.8	125	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	131	0.00
8	S	Nitrobenzene-d5	0.140	0.142	-1.4	127	0.00
9	S	2-Nitrophenol-d4	0.150	0.162	-8.0	138	0.00
10	S	2,4-Dichlorophenol-d3	0.265	0.282	-6.4	136	0.00
11		Naphthalene	0.973	0.967	0.6	127	0.00
12	S	4-Chloroaniline-d4	0.208	0.306	-47.1#	207#	0.00
13		2-Methylnaphthalene	0.705	0.706	-0.1	129	0.00
14		1-Methylnaphthalene	0.721	0.711	1.4	124	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	125	0.00
16	S	Dimethylphthalate-d6	1.542	1.477	4.2	119	0.00
17	S	Acenaphthylene-d8	1.986	2.017	-1.6	123	0.00
18		Acenaphthylene	2.068	2.014	2.6	119	0.00
19	C	Acenaphthene	1.420	1.387	2.3	120	0.00
20	S	4-Nitrophenol-d4	0.271	0.241	11.1	116	0.00
21	S	Fluorene-d10	1.388	1.361	1.9	122	0.00
22		Fluorene	1.573	1.496	4.9	117	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	116	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.099	0.112	-13.1	141	0.00
25		Phenanthrene	1.069	1.051	1.7	112	0.00
26	S	Anthracene-d10	0.921	0.904	1.8	113	0.00
27		Anthracene	1.074	1.072	0.2	112	0.00
28	C	Fluoranthene	1.111	1.060	4.6	108	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	111	0.00
30	S	Pyrene-d10	1.055	1.058	-0.3	109	0.00
31		Pyrene	1.377	1.337	2.9	105	0.00
32		Benzo(a)anthracene	1.109	1.086	2.1	106	0.00
33		Chrysene	1.079	1.020	5.5	104	0.00
34	I	Perylene-d12	1.000	1.000	0.0	116	-0.02
35		Benzo(b)fluoranthene	1.150	1.068	7.1	106	-0.02
36		Benzo(k)fluoranthene	1.131	1.064	5.9	108	0.00
37	S	Benzo(a)pyrene-d12	0.888	0.857	3.5	111	0.00
38	C	Benzo(a)pyrene	1.099	1.036	5.7	108	0.00
39		Indeno(1,2,3-cd)pyrene	1.250	1.223	2.2	114	-0.02
40		Dibenzo(a,h)anthracene	1.035	1.031	0.4	114	-0.02
41		Benzo(g,h,i)perylene	1.057	0.993	6.1	107	-0.02

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