Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG102016\
Data File : BG024472.D
Acq On : 20 Oct 2016 21:14
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : SSTD02014

Quant Time: Oct 21 17:47:43 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\SOM02.2-EPA-BG102016-MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Fri Oct 21 17:27:43 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	0.374	0.361	3.5	97	0.00
3	S	Phenol-d5	1.819	1.851	-1.8	99	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	1.124	1.124	0.0	93	0.00
5	S	2-Chlorophenol-d4	1.280	1.311	-2.4	98	0.00
6	S	4-Methylphenol-d8	1.517	1.574	-3.8	100	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
8	S	Nitrobenzene-d5	0.150	0.151	-0.7	104	0.00
9	S	2-Nitrophenol-d4	0.174	0.173	0.6	96	0.00
10	S	2,4-Dichlorophenol-d3	0.353	0.354	-0.3	98	0.00
11	S	4-Chloroaniline-d4	0.367	0.414	-12.8	106	0.00
12		1-Methylnaphthalene	0.743	0.761	-2.4	101	0.00
13	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
14	S	Dimethylphthalate-d6	1.349	1.372	-1.7	103	0.00
15	S	Acenaphthylene-d8	1.605	1.643	-2.4	100	0.00
16	S	4-Nitrophenol-d4	0.231	0.232	-0.4	106	0.00
17	S	Fluorene-d10	1.285	1.326	-3.2	107	0.00
18	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
19	S	4,6-Dinitro-2-methylphenol-	0.134	0.131	2.2	108	0.00
20	S	Anthracene-d10	0.889	0.916	-3.0	103	0.00
21		1,2,3,4-Tetrachlorobenzene	0.305	0.304	0.3	98	0.00
22		Pentachlorobenzene	0.320	0.323	-0.9	102	0.00
23	I	Chrysene-d12	1.000	1.000	0.0	99	0.00
24	S	Pyrene-d10	0.869	0.929	-6.9	103	0.00
25	I	Perylene-d12	1.000	1.000	0.0	103	0.00
26	S	Benzo(a)pyrene-d12	0.875	0.880	-0.6	104	0.00

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