Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM061016\
Data File : BM005718.D
Acq On : 11 Jun 2016 00:04
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : SSTDCCC020

Quant Time: Jun 13 11:55:35 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM061016.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jun 13 03:29:45 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	104	0.00
2	S	1,4-Dioxane-d8	0.461	0.444	3.7	103	0.00
3	S	Phenol-d5	1.726	1.697	1.7	102	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	1.037	1.051	-1.4	103	0.00
5	S	2-Chlorophenol-d4	1.395	1.408	-0.9	103	0.00
6	S	4-Methylphenol-d8	1.458	1.469	-0.8	104	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
8	S	Nitrobenzene-d5	0.144	0.145	-0.7	103	0.00
9	S	2-Nitrophenol-d4	0.163	0.165	-1.2	100	0.00
10	S	2,4-Dichlorophenol-d3	0.296	0.299	-1.0	105	0.00
11		Naphthalene	1.002	0.982	2.0	103	0.00
12	S	4-Chloroaniline-d4	0.357	0.426	-19.3	104	0.00
13		2-Methylnaphthalene	0.754	0.743	1.5	103	0.00
14		1-Methylnaphthalene	0.745	0.736	1.2	103	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
16	S	Dimethylphthalate-d6	1.650	1.628	1.3	103	0.00
17	S	Acenaphthylene-d8	2.018	2.017	0.0	103	0.00
18		Acenaphthylene	2.066	2.050	0.8	103	0.00
19	C	Acenaphthene	1.344	1.332	0.9	103	0.00
20	S	4-Nitrophenol-d4	0.241	0.242	-0.4	104	-0.01
21	S	Fluorene-d10	1.397	1.390	0.5	103	-0.01
22		Fluorene	1.534	1.539	-0.3	103	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	104	-0.01
24	S	4,6-Dinitro-2-methylphenol-	0.103	0.099	3.9	101	-0.02
25		Phenanthrene	1.095	1.089	0.5	102	-0.01
26	S	Anthracene-d10	0.942	0.944	-0.2	103	0.00
27		Anthracene	1.111	1.136	-2.3	103	0.00
28	C	Fluoranthene	1.286	1.291	-0.4	102	-0.01
29	I	Chrysene-d12	1.000	1.000	0.0	103	-0.01
30	S	Pyrene-d10	0.961	0.928	3.4	103	0.00
31		Pyrene	1.217	1.191	2.1	103	0.00
32		Benzo(a)anthracene	1.155	1.144	1.0	103	0.00
33		Chrysene	1.139	1.138	0.1	102	0.00
34	I	Perylene-d12	1.000	1.000	0.0	103	-0.01
35		Benzo(b)fluoranthene	1.188	1.192	-0.3	106	-0.02
36		Benzo(k)fluoranthene	1.201	1.186	1.2	100	-0.01
37	S	Benzo(a)pyrene-d12	0.912	0.913	-0.1	103	-0.02
38	C	Benzo(a)pyrene	1.128	1.151	-2.0	105	-0.02
39		Indeno(1,2,3-cd)pyrene	1.101	1.113	-1.1	104	-0.03
40		Dibenzo(a,h)anthracene	0.921	0.931	-1.1	105	-0.02
41		Benzo(g,h,i)perylene	0.921	0.904	1.8	104	-0.02

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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