Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM062616\
Data File : BM006033.D
Acq On : 26 Jun 2016 10:48
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02035

Quant Time: Jun 26 11:18:55 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM062516.M
Quant Title : SVOA CALIBRATION
QLast Update : Sun Jun 26 04:58:10 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	71	0.00
2	S	1,4-Dioxane-d8	0.448	0.437	2.5	67	0.00
3	S	Phenol-d5	2.104	1.981	5.8	67	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	1.170	1.153	1.5	68	0.00
5	S	2-Chlorophenol-d4	1.465	1.467	-0.1	70	0.00
6	S	4-Methylphenol-d8	1.776	1.634	8.0	66	-0.01
7	I	Naphthalene-d8	1.000	1.000	0.0	70	0.00
8	S	Nitrobenzene-d5	0.152	0.143	5.9	65	0.00
9	S	2-Nitrophenol-d4	0.169	0.150	11.2	61	0.00
10	S	2,4-Dichlorophenol-d3	0.325	0.323	0.6	68	0.00
11		Naphthalene	0.999	1.009	-1.0	69	0.00
12	S	4-Chloroaniline-d4	0.353	0.423	-19.8	67	0.00
13		2-Methylnaphthalene	0.804	0.814	-1.2	69	0.00
14		1-Methylnaphthalene	0.770	0.779	-1.2	69	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	68	0.00
16	S	Dimethylphthalate-d6	1.712	1.651	3.6	65	0.00
17	S	Acenaphthylene-d8	1.986	2.006	-1.0	67	0.00
18		Acenaphthylene	2.094	2.150	-2.7	68	0.00
19	C	Acenaphthene	1.330	1.370	-3.0	69	0.00
20	S	4-Nitrophenol-d4	0.365	0.396	-8.5	69	-0.02
21	S	Fluorene-d10	1.462	1.505	-2.9	68	0.00
22		Fluorene	1.602	1.692	-5.6	69	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	72	-0.01
24	S	4,6-Dinitro-2-methylphenol-	0.111	0.094	15.3	59	-0.02
25		Phenanthrene	1.076	1.114	-3.5	73	-0.01
26	S	Anthracene-d10	0.919	0.943	-2.6	72	0.00
27		Anthracene	1.111	1.149	-3.4	73	-0.01
28	C	Fluoranthene	1.378	1.551	-12.6	78	-0.01
29	I	Chrysene-d12	1.000	1.000	0.0	91	-0.01
30	S	Pyrene-d10	0.876	0.797	9.0	80	-0.01
31		Pyrene	1.145	1.058	7.6	81	-0.01
32		Benzo(a)anthracene	1.188	1.208	-1.7	89	-0.01
33		Chrysene	1.085	1.130	-4.1	91	-0.01
34	I	Perylene-d12	1.000	1.000	0.0	97	-0.01
35		Benzo(b)fluoranthene	1.162	1.170	-0.7	96	-0.02
36		Benzo(k)fluoranthene	1.081	1.091	-0.9	94	-0.02
37	S	Benzo(a)pyrene-d12	0.903	0.919	-1.8	97	-0.02
38	C	Benzo(a)pyrene	1.114	1.144	-2.7	98	-0.02
39		Indeno(1,2,3-cd)pyrene	1.328	1.382	-4.1	97	-0.04
40		Dibenzo(a,h)anthracene	1.104	1.149	-4.1	97	-0.02
41		Benzo(g,h,i)perylene	1.149	1.186	-3.2	97	-0.02

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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