Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\Data\BM072016\
Data File : BM006561.D
Acq On : 20 Jul 2016 19:09
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02054

Quant Time: Jul 21 01:01:58 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jul 20 18:59:53 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	0.495	0.506	-2.2	106	0.00
3	S	Phenol-d5	1.702	1.721	-1.1	105	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	1.054	1.090	-3.4	105	0.00
5	S	2-Chlorophenol-d4	1.360	1.363	-0.2	105	0.00
6	S	4-Methylphenol-d8	1.320	1.350	-2.3	105	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	105	0.00
8	S	Nitrobenzene-d5	0.151	0.152	-0.7	107	0.00
9	S	2-Nitrophenol-d4	0.161	0.164	-1.9	107	0.00
10	S	2,4-Dichlorophenol-d3	0.304	0.306	-0.7	105	0.00
11		Naphthalene	1.031	1.021	1.0	105	0.00
12	S	4-Chloroaniline-d4	0.335	0.409	-22.1#	104	0.00
13		2-Methylnaphthalene	0.757	0.741	2.1	102	0.00
14		1-Methylnaphthalene	0.722	0.710	1.7	103	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
16	S	Dimethylphthalate-d6	1.568	1.583	-1.0	102	0.00
17	S	Acenaphthylene-d8	2.002	2.049	-2.3	103	0.00
18		Acenaphthylene	2.144	2.161	-0.8	101	0.00
19	C	Acenaphthene	1.378	1.385	-0.5	101	0.00
20	S	4-Nitrophenol-d4	0.272	0.285	-4.8	107	0.00
21	S	Fluorene-d10	1.402	1.413	-0.8	102	0.00
22		Fluorene	1.552	1.600	-3.1	101	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.104	0.101	2.9	98	0.00
25		Phenanthrene	1.125	1.142	-1.5	101	0.00
26	S	Anthracene-d10	0.950	0.966	-1.7	100	0.00
27		Anthracene	1.158	1.197	-3.4	103	0.00
28	C	Fluoranthene	1.321	1.376	-4.2	102	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
30	S	Pyrene-d10	0.911	0.910	0.1	102	0.00
31		Pyrene	1.211	1.214	-0.2	101	0.00
32		Benzo(a)anthracene	1.212	1.216	-0.3	100	0.00
33		Chrysene	1.118	1.129	-1.0	101	0.00
34	I	Perylene-d12	1.000	1.000	0.0	100	0.00
35		Benzo(b)fluoranthene	1.199	1.214	-1.3	101	0.00
36		Benzo(k)fluoranthene	1.159	1.185	-2.2	100	0.00
37	S	Benzo(a)pyrene-d12	0.920	0.938	-2.0	101	0.00
38	C	Benzo(a)pyrene	1.167	1.191	-2.1	100	0.00
39		Indeno(1,2,3-cd)pyrene	1.351	1.386	-2.6	101	0.00
40		Dibenzo(a,h)anthracene	1.120	1.159	-3.5	101	0.00
41		Benzo(g,h,i)perylene	1.158	1.182	-2.1	101	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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