Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM073116\
Data File : BM006813.D
Acq On : 01 Aug 2016 06:12
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02064

Quant Time: Aug 01 06:47:04 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 30 01:37:26 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	155#	0.00
2	S	1,4-Dioxane-d8	0.495	0.452	8.7	140	0.00
3	S	Phenol-d5	1.702	1.788	-5.1	161#	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	1.054	1.099	-4.3	157#	0.00
5	S	2-Chlorophenol-d4	1.360	1.394	-2.5	160#	0.00
6	S	4-Methylphenol-d8	1.320	1.390	-5.3	160#	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	158#	0.00
8	S	Nitrobenzene-d5	0.151	0.161	-6.6	171#	0.00
9	S	2-Nitrophenol-d4	0.161	0.174	-8.1	171#	0.00
10	S	2,4-Dichlorophenol-d3	0.304	0.315	-3.6	162#	0.00
11		Naphthalene	1.031	1.013	1.7	156#	0.00
12	S	4-Chloroaniline-d4	0.335	0.388	-15.8	148	0.00
13		2-Methylnaphthalene	0.757	0.738	2.5	153#	0.00
14		1-Methylnaphthalene	0.722	0.701	2.9	153#	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	151#	0.00
16	S	Dimethylphthalate-d6	1.568	1.579	-0.7	152#	0.00
17	S	Acenaphthylene-d8	2.002	2.068	-3.3	156#	0.00
18		Acenaphthylene	2.144	2.155	-0.5	151#	0.00
19	C	Acenaphthene	1.378	1.372	0.4	150	0.00
20	S	4-Nitrophenol-d4	0.272	0.274	-0.7	154#	0.01
21	S	Fluorene-d10	1.402	1.391	0.8	150	0.00
22		Fluorene	1.552	1.551	0.1	147	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	145	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.104	0.098	5.8	137	0.01
25		Phenanthrene	1.125	1.117	0.7	142	0.00
26	S	Anthracene-d10	0.950	0.965	-1.6	143	0.00
27		Anthracene	1.158	1.162	-0.3	143	0.00
28	C	Fluoranthene	1.321	1.310	0.8	139	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	127	0.00
30	S	Pyrene-d10	0.911	0.986	-8.2	139	0.00
31		Pyrene	1.211	1.289	-6.4	135	0.00
32		Benzo(a)anthracene	1.212	1.212	0.0	125	0.00
33		Chrysene	1.118	1.106	1.1	125	0.00
34	I	Perylene-d12	1.000	1.000	0.0	128	0.01
35		Benzo(b)fluoranthene	1.199	1.195	0.3	126	0.00
36		Benzo(k)fluoranthene	1.159	1.132	2.3	122	0.01
37	S	Benzo(a)pyrene-d12	0.920	0.925	-0.5	127	0.00
38	C	Benzo(a)pyrene	1.167	1.165	0.2	125	0.01
39		Indeno(1,2,3-cd)pyrene	1.351	1.362	-0.8	126	0.02
40		Dibenzo(a,h)anthracene	1.120	1.138	-1.6	126	0.02
41		Benzo(g,h,i)perylene	1.158	1.186	-2.4	129	0.02

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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