Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM060716\
Data File : BM005690.D
Acq On : 07 Jun 2016 13:41
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02048

Quant Time: Jun 07 14:13:33 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM060216.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Jun 07 04:25:20 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	101	0.00
2	S	1,4-Dioxane-d8	0.476	0.449	5.7	98	0.00
3	S	Phenol-d5	1.758	1.892	-7.6	108	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	1.070	1.123	-5.0	104	0.00
5	S	2-Chlorophenol-d4	1.451	1.499	-3.3	105	0.00
6	S	4-Methylphenol-d8	1.391	1.581	-13.7	115	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
8	S	Nitrobenzene-d5	0.154	0.157	-1.9	115	0.00
9	S	2-Nitrophenol-d4	0.164	0.174	-6.1	117	0.00
10	S	2,4-Dichlorophenol-d3	0.292	0.309	-5.8	119	0.00
11		Naphthalene	1.017	1.011	0.6	112	0.00
12	S	4-Chloroaniline-d4	0.347	0.446	-28.5#	129	0.00
13		2-Methylnaphthalene	0.707	0.741	-4.8	118	0.00
14		1-Methylnaphthalene	0.695	0.737	-6.0	119	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	130	0.00
16	S	Dimethylphthalate-d6	1.538	1.641	-6.7	141	0.00
17	S	Acenaphthylene-d8	2.000	2.018	-0.9	132	0.00
18		Acenaphthylene	2.069	2.089	-1.0	131	0.00
19	C	Acenaphthene	1.343	1.355	-0.9	132	0.00
20	S	4-Nitrophenol-d4	0.220	0.276	-25.5#	167#	0.00
21	S	Fluorene-d10	1.321	1.409	-6.7	140	0.00
22		Fluorene	1.453	1.545	-6.3	137	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	154#	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.094	0.097	-3.2	173#	0.00
25		Phenanthrene	1.104	1.089	1.4	152#	0.00
26	S	Anthracene-d10	0.950	0.952	-0.2	153#	0.00
27		Anthracene	1.122	1.122	0.0	153#	0.00
28	C	Fluoranthene	1.128	1.259	-11.6	170#	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	173#	0.00
30	S	Pyrene-d10	1.005	0.963	4.2	167#	0.00
31		Pyrene	1.262	1.211	4.0	168#	0.00
32		Benzo(a)anthracene	1.173	1.194	-1.8	175#	0.00
33		Chrysene	1.108	1.115	-0.6	176#	0.00
34	I	Perylene-d12	1.000	1.000	0.0	175#	0.00
35		Benzo(b)fluoranthene	1.206	1.250	-3.6	186#	0.00
36		Benzo(k)fluoranthene	1.132	1.106	2.3	170#	0.00
37	S	Benzo(a)pyrene-d12	0.920	0.929	-1.0	179#	0.00
38	C	Benzo(a)pyrene	1.161	1.147	1.2	173#	0.00
39		Indeno(1,2,3-cd)pyrene	1.374	1.189	13.5	152#	0.00
40		Dibenzo(a,h)anthracene	1.153	1.000	13.3	151#	0.00
41		Benzo(g,h,i)perylene	1.152	0.951	17.4	147	0.00

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