Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM080616\
Data File : BM006943.D
Acq On : 06 Aug 2016 18:21
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02022

Quant Time: Aug 08 03:51:33 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM080616.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Aug 08 03:51:46 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	107	0.00
2	S	1,4-Dioxane-d8	0.438	0.445	-1.6	104	0.00
3	S	Phenol-d5	1.896	1.911	-0.8	113	0.00
4	S	Bis-(2-Chloroethyl)ether-d8	1.189	1.229	-3.4	110	0.00
5	S	2-Chlorophenol-d4	1.354	1.438	-6.2	112	0.00
6	S	4-Methylphenol-d8	1.583	1.606	-1.5	113	0.00
7	I	Naphthalene-d8	1.000	1.000	0.0	116	0.00
8	S	Nitrobenzene-d5	0.149	0.151	-1.3	115	0.00
9	S	2-Nitrophenol-d4	0.178	0.182	-2.2	112	0.00
10	S	2,4-Dichlorophenol-d3	0.337	0.350	-3.9	118	0.00
11		Naphthalene	1.027	1.029	-0.2	114	0.00
12	S	4-Chloroaniline-d4	0.403	0.437	-8.4	116	0.00
13		2-Methylnaphthalene	0.822	0.840	-2.2	116	0.00
14		1-Methylnaphthalene	0.790	0.799	-1.1	115	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	114	0.00
16	S	Dimethylphthalate-d6	1.760	1.783	-1.3	114	0.00
17	S	Acenaphthylene-d8	2.036	2.087	-2.5	115	0.00
18		Acenaphthylene	2.136	2.181	-2.1	115	0.00
19	C	Acenaphthene	1.393	1.411	-1.3	115	0.00
20	S	4-Nitrophenol-d4	0.255	0.208	18.4	112	0.00
21	S	Fluorene-d10	1.496	1.502	-0.4	115	0.00
22		Fluorene	1.761	1.763	-0.1	114	0.00
23	I	Phenanthrene-d10	1.000	1.000	0.0	111	0.00
24	S	4,6-Dinitro-2-methylphenol-	0.083	0.076	8.4	140	0.00
25		Phenanthrene	1.119	1.124	-0.4	111	0.00
26	S	Anthracene-d10	0.966	0.960	0.6	110	0.00
27		Anthracene	1.171	1.172	-0.1	111	0.00
28	C	Fluoranthene	1.361	1.350	0.8	107	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	91	0.00
30	S	Pyrene-d10	0.953	1.028	-7.9	103	0.00
31		Pyrene	1.243	1.336	-7.5	103	0.00
32		Benzo(a)anthracene	1.187	1.196	-0.8	92	0.00
33		Chrysene	1.045	1.053	-0.8	91	0.00
34	I	Perylene-d12	1.000	1.000	0.0	77	0.00
35		Benzo(b)fluoranthene	1.305	1.325	-1.5	81	0.00
36		Benzo(k)fluoranthene	1.260	1.224	2.9	74	0.00
37	S	Benzo(a)pyrene-d12	0.953	0.956	-0.3	76	0.00
38	C	Benzo(a)pyrene	1.184	1.178	0.5	77	0.00
39		Indeno(1,2,3-cd)pyrene	1.201	1.284	-6.9	83	0.00
40		Dibenzo(a,h)anthracene	1.011	1.071	-5.9	82	0.00
41		Benzo(g,h,i)perylene	0.946	1.040	-9.9	87	-0.01

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