Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM073116\
Data File : BM006825.D
Acq On : 01 Aug 2016 14:05
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02065

Quant Time: Aug 01 16:00:44 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	207#	0.00
2	S	1,4-Dioxane-d8	0.495	0.451	8.9	187#	0.00
3	S	Phenol-d5	1.702	1.767	-3.8	213#	0.01
4	S	Bis-(2-Chloroethyl)ether-d8	1.054	1.075	-2.0	205#	0.00
5	S	2-Chlorophenol-d4	1.360	1.380	-1.5	211#	0.00
6	S	4-Methylphenol-d8	1.320	1.351	-2.3	208#	0.01
7	I	Naphthalene-d8	1.000	1.000	0.0	207#	0.00
8	S	Nitrobenzene-d5	0.151	0.161	-6.6	225#	0.00
9	S	2-Nitrophenol-d4	0.161	0.174	-8.1	225#	0.00
10	S	2,4-Dichlorophenol-d3	0.304	0.316	-3.9	214#	0.00
11		Naphthalene	1.031	1.011	1.9	206#	0.00
12	S	4-Chloroaniline-d4	0.335	0.395	-17.9	199#	0.00
13		2-Methylnaphthalene	0.757	0.731	3.4	200#	0.00
14		1-Methylnaphthalene	0.722	0.687	4.8	197#	0.00
15	I	Acenaphthene-d10	1.000	1.000	0.0	195#	0.00
16	S	Dimethylphthalate-d6	1.568	1.569	-0.1	195#	0.01
17	S	Acenaphthylene-d8	2.002	2.063	-3.0	200#	0.00
18		Acenaphthylene	2.144	2.149	-0.2	194#	0.00
19	C	Acenaphthene	1.378	1.355	1.7	190#	0.00
20	S	4-Nitrophenol-d4	0.272	0.266	2.2	193#	0.02
21	S	Fluorene-d10	1.402	1.383	1.4	192#	0.00
22		Fluorene	1.552	1.528	1.5	186#	0.01
23	I	Phenanthrene-d10	1.000	1.000	0.0	183#	0.01
24	S	4,6-Dinitro-2-methylphenol-	0.104	0.064	38.5#	113	0.02
25		Phenanthrene	1.125	1.117	0.7	180#	0.01
26	S	Anthracene-d10	0.950	0.960	-1.1	181#	0.01
27		Anthracene	1.158	1.157	0.1	180#	0.00
28	C	Fluoranthene	1.321	1.278	3.3	172#	0.00
29	I	Chrysene-d12	1.000	1.000	0.0	150	0.01
30	S	Pyrene-d10	0.911	1.026	-12.6	170#	0.00
31		Pyrene	1.211	1.343	-10.9	165#	0.01
32		Benzo(a)anthracene	1.212	1.215	-0.2	147	0.01
33		Chrysene	1.118	1.139	-1.9	151#	0.01
34	I	Perylene-d12	1.000	1.000	0.0	155#	0.02
35		Benzo(b)fluoranthene	1.199	1.156	3.6	148	0.02
36		Benzo(k)fluoranthene	1.159	1.136	2.0	148	0.02
37	S	Benzo(a)pyrene-d12	0.920	0.916	0.4	152#	0.02
38	C	Benzo(a)pyrene	1.167	1.148	1.6	150	0.02
39		Indeno(1,2,3-cd)pyrene	1.351	1.350	0.1	152#	0.04
40		Dibenzo(a,h)anthracene	1.120	1.123	-0.3	151#	0.03
41		Benzo(g,h,i)perylene	1.158	1.174	-1.4	155#	0.04

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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