Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072216\
Data File : BM006643.D
Acq On : 23 Jul 2016 02:46
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02063

Quant Time: Jul 23 03:28:27 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 23 00:16:38 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	265#	0.01
2	S	1,4-Dioxane-d8	0.495	0.466	5.9	247#	0.00
3	S	Phenol-d5	1.702	1.787	-5.0	276#	0.01
4	S	Bis-(2-Chloroethyl)ether-d8	1.054	1.098	-4.2	269#	0.01
5	S	2-Chlorophenol-d4	1.360	1.403	-3.2	275#	0.00
6	S	4-Methylphenol-d8	1.320	1.363	-3.3	269#	0.02
7	I	Naphthalene-d8	1.000	1.000	0.0	261#	0.01
8	S	Nitrobenzene-d5	0.151	0.163	-7.9	285#	0.01
9	S	2-Nitrophenol-d4	0.161	0.173	-7.5	281#	0.00
10	S	2,4-Dichlorophenol-d3	0.304	0.316	-3.9	270#	0.01
11		Naphthalene	1.031	1.022	0.9	262#	0.01
12	S	4-Chloroaniline-d4	0.335	0.414	-23.6#	262#	0.01
13		2-Methylnaphthalene	0.757	0.736	2.8	254#	0.01
14		1-Methylnaphthalene	0.722	0.707	2.1	256#	0.01
15	I	Acenaphthene-d10	1.000	1.000	0.0	245#	0.00
16	S	Dimethylphthalate-d6	1.568	1.570	-0.1	245#	0.00
17	S	Acenaphthylene-d8	2.002	2.064	-3.1	252#	0.01
18		Acenaphthylene	2.144	2.151	-0.3	244#	0.01
19	C	Acenaphthene	1.378	1.372	0.4	243#	0.01
20	S	4-Nitrophenol-d4	0.272	0.282	-3.7	257#	0.02
21	S	Fluorene-d10	1.402	1.397	0.4	244#	0.01
22		Fluorene	1.552	1.524	1.8	234#	0.01
23	I	Phenanthrene-d10	1.000	1.000	0.0	236#	0.01
24	S	4,6-Dinitro-2-methylphenol-	0.104	0.065	37.5#	148	0.02
25		Phenanthrene	1.125	1.119	0.5	231#	0.01
26	S	Anthracene-d10	0.950	0.953	-0.3	231#	0.01
27		Anthracene	1.158	1.145	1.1	229#	0.01
28	C	Fluoranthene	1.321	1.328	-0.5	230#	0.01
29	I	Chrysene-d12	1.000	1.000	0.0	203#	0.01
30	S	Pyrene-d10	0.911	1.013	-11.2	228#	0.01
31		Pyrene	1.211	1.322	-9.2	221#	0.01
32		Benzo(a)anthracene	1.212	1.244	-2.6	205#	0.01
33		Chrysene	1.118	1.176	-5.2	211#	0.01
34	I	Perylene-d12	1.000	1.000	0.0	209#	0.02
35		Benzo(b)fluoranthene	1.199	1.202	-0.3	207#	0.02
36		Benzo(k)fluoranthene	1.159	1.200	-3.5	211#	0.01
37	S	Benzo(a)pyrene-d12	0.920	0.958	-4.1	214#	0.02
38	C	Benzo(a)pyrene	1.167	1.202	-3.0	211#	0.02
39		Indeno(1,2,3-cd)pyrene	1.351	1.398	-3.5	212#	0.04
40		Dibenzo(a,h)anthracene	1.120	1.166	-4.1	211#	0.04
41		Benzo(g,h,i)perylene	1.158	1.231	-6.3	220#	0.04

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM072216\
Data File : BM006643.D
Acq On : 23 Jul 2016 02:46
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02063

Quant Time: Jul 23 03:28:27 2016
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-BM072016.M
Quant Title : SVOA CALIBRATION
QLast Update : Sat Jul 23 00:16:38 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)

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