Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092616\
Data File : BM007592.D
Acq On : 27 Sep 2016 03:43
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02056

Quant Time: Sep 27 04:16:53 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092316.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 27 03:37:16 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	135	0.00
2		1,4-Dioxane	0.427	0.471	-10.3	138	0.00
3	S	1,4-Dioxane-d8	0.407	0.431	-5.9	135	0.00
4		Benzaldehyde	1.191	1.235	-3.7	135	0.00
5	S	Phenol-d5	1.775	1.770	0.3	138	0.00
6		Phenol	1.801	1.807	-0.3	137	0.00
7	S	Bis-(2-Chloroethyl)ether-d8	0.954	0.997	-4.5	137	0.00
8		Bis(2-Chloroethyl)ether	1.290	1.338	-3.7	137	0.00
9	S	2-Chlorophenol-d4	1.289	1.312	-1.8	132	0.00
10		2-Chlorophenol	1.293	1.338	-3.5	134	0.00
11		2-Methylphenol	1.388	1.371	1.2	135	0.00
12		2,2'-oxybis(1-Chloropropane	1.403	1.476	-5.2	135	0.00
13	S	4-Methylphenol-d8	1.515	1.511	0.3	135	0.00
14		Acetophenone	2.308	2.354	-2.0	133	0.00
15	P	N-Nitroso-di-n-propylamine	1.172	1.223	-4.4	133	0.00
16		4-Methylphenol	1.554	1.555	-0.1	134	0.00
17		Hexachloroethane	0.529	0.536	-1.3	134	0.00
18	I	Naphthalene-d8	1.000	1.000	0.0	135	0.00
19	S	Nitrobenzene-d5	0.143	0.145	-1.4	132	0.00
20		Nitrobenzene	0.366	0.382	-4.4	136	0.00
21		Isophorone	0.701	0.718	-2.4	133	0.00
22	S	2-Nitrophenol-d4	0.162	0.164	-1.2	133	0.00
23	C	2-Nitrophenol	0.169	0.172	-1.8	131	0.00
24		2,4-Dimethylphenol	0.391	0.401	-2.6	133	0.00
25		Bis(2-Chloroethoxy)methane	0.400	0.408	-2.0	133	0.00
26	S	2,4-Dichlorophenol-d3	0.324	0.335	-3.4	135	0.00
27	C	2,4-Dichlorophenol	0.319	0.328	-2.8	133	0.00
28		Naphthalene	0.966	0.989	-2.4	135	0.00
29	S	4-Chloroaniline-d4	0.340	0.380	-11.8	149	0.00
30		4-Chloroaniline	0.356	0.388	-9.0	144	0.00
31	C	Hexachlorobutadiene	0.233	0.228	2.1	129	0.00
32		Caprolactam	0.117	0.112	4.3	132	0.00
33	C	4-Chloro-3-methylphenol	0.388	0.400	-3.1	134	0.00
34		2-Methylnaphthalene	0.758	0.777	-2.5	134	0.00
35	I	Acenaphthene-d10	1.000	1.000	0.0	133	0.00
36		1,2,4,5-Tetrachlorobenzene	0.590	0.602	-2.0	134	0.00
37		Hexachlorocyclopentadiene	0.317	0.265	16.4	120	0.00
38	C	2,4,6-Trichlorophenol	0.376	0.387	-2.9	132	0.00
39		2,4,5-Trichlorophenol	0.412	0.431	-4.6	130	0.00
40		1,1'-Biphenyl	1.328	1.394	-5.0	135	0.00
41		2-Chloronaphthalene	1.033	1.062	-2.8	132	0.00
42		2-Nitroaniline	0.318	0.345	-8.5	136	0.00
43	S	Dimethylphthalate-d6	1.455	1.482	-1.9	130	0.00
44		Dimethylphthalate	1.411	1.426	-1.1	130	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092616\
Data File : BM007592.D
Acq On : 27 Sep 2016 03:43
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02056

Quant Time: Sep 27 04:16:53 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092316.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 27 03:37:16 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)
45		2,6-Dinitrotoluene	0.280	0.289	-3.2	131	0.00
46	S	Acenaphthylene-d8	1.688	1.747	-3.5	132	0.00
47		Acenaphthylene	1.676	1.741	-3.9	132	0.00
48		3-Nitroaniline	0.284	0.301	-6.0	137	0.00
49	C	Acenaphthene	1.152	1.183	-2.7	132	0.00
50		2,4-Dinitrophenol	0.187	0.154	17.6	120	0.00
51	S	4-Nitrophenol-d4	0.257	0.255	0.8	133	0.00
52		4-Nitrophenol	0.267	0.265	0.7	129	0.00
53		Dibenzofuran	1.700	1.759	-3.5	133	0.00
54		2,4-Dinitrotoluene	0.432	0.452	-4.6	131	0.00
55		2,3,4,6-Tetrachlorophenol	0.407	0.427	-4.9	133	0.00
56		Diethylphthalate	1.440	1.471	-2.2	131	0.00
57	S	Fluorene-d10	1.276	1.319	-3.4	132	0.00
58		Fluorene	1.403	1.442	-2.8	131	0.00
59		4-Chlorophenyl-phenylether	0.738	0.756	-2.4	132	0.00
60		4-Nitroaniline	0.313	0.333	-6.4	139	0.00
61	I	Phenanthrene-d10	1.000	1.000	0.0	135	0.00
62	S	4,6-Dinitro-2-methylphenol-	0.122	0.113	7.4	132	0.00
63		4,6-Dinitro-2-methylphenol	0.125	0.117	6.4	131	0.00
64		N-Nitrosodiphenylamine	0.540	0.550	-1.9	133	0.00
65		4-Bromophenyl-phenylether	0.218	0.219	-0.5	130	0.00
66		Hexachlorobenzene	0.246	0.245	0.4	131	0.00
67		Atrazine	0.225	0.227	-0.9	133	0.00
68	C	Pentachlorophenol	0.154	0.146	5.2	133	0.00
69		Phenanthrene	1.059	1.080	-2.0	133	0.00
70	S	Anthracene-d10	0.919	0.944	-2.7	134	0.00
71		Anthracene	1.075	1.107	-3.0	134	0.00
72		Carbazole	0.956	1.007	-5.3	135	0.00
73		Di-n-butylphthalate	1.081	1.117	-3.3	132	0.00
74	C	Fluoranthene	1.336	1.428	-6.9	133	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	131	0.00
76	S	Pyrene-d10	0.794	0.832	-4.8	134	0.00
77		Pyrene	0.981	1.027	-4.7	134	0.00
78		Butylbenzylphthalate	0.377	0.398	-5.6	132	0.00
79		3,3'-Dichlorobenzidine	0.367	0.375	-2.2	136	0.00
80		Benzo(a)anthracene	1.093	1.125	-2.9	131	0.00
81		Bis(2-ethylhexyl)phthalate	0.543	0.578	-6.4	131	0.00
82		Chrysene	1.044	1.077	-3.2	131	0.00
83	I	Perylene-d12	1.000	1.000	0.0	123	0.00
84		Di-n-octyl phthalate	0.966	1.080	-11.8	131	0.00
85		Benzo(b)fluoranthene	1.122	1.169	-4.2	121	0.00
86		Benzo(k)fluoranthene	1.079	1.159	-7.4	132	0.00
87	S	Benzo(a)pyrene-d12	0.885	0.916	-3.5	123	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM092616\
Data File : BM007592.D
Acq On : 27 Sep 2016 03:43
Operator : UM/SJ
Sample : SSTDCCC020EC
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02056

Quant Time: Sep 27 04:16:53 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM092316.M
Quant Title : SVOA CALIBRATION
QLast Update : Tue Sep 27 03:37:16 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)
88 C	Benzo(a)pyrene	1.077	1.120	-4.0	123	0.00
89	Indeno(1,2,3-cd)pyrene	1.189	1.181	0.7	116	0.00
90	Dibenzo(a,h)anthracene	0.980	0.973	0.7	116	0.00
91	Benzo(g,h,i)perylene	0.998	0.991	0.7	117	0.00

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