Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM081816\
Data File : BM007176.D
Acq On : 17 Aug 2016 12:39
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02056

Quant Time: Aug 17 13:54:30 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM081116.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 17 04:01:43 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	114	0.00
2		1,4-Dioxane	0.476	0.473	0.6	116	-0.01
3	S	1,4-Dioxane-d8	0.458	0.460	-0.4	113	0.00
4		Benzaldehyde	0.974	1.039	-6.7	113	0.00
5	S	Phenol-d5	1.620	1.594	1.6	114	0.00
6		Phenol	1.707	1.660	2.8	112	0.00
7	S	Bis-(2-Chloroethyl)ether-d8	0.929	0.922	0.8	112	0.00
8		Bis(2-Chloroethyl)ether	1.259	1.252	0.6	112	0.00
9	S	2-Chlorophenol-d4	1.290	1.301	-0.9	115	0.00
10		2-Chlorophenol	1.320	1.328	-0.6	113	0.00
11		2-Methylphenol	1.283	1.231	4.1	112	0.00
12		2,2'-oxybis(1-Chloropropane	1.443	1.387	3.9	111	0.00
13	S	4-Methylphenol-d8	1.333	1.260	5.5	111	0.00
14		Acetophenone	2.053	2.013	1.9	112	0.00
15	P	N-Nitroso-di-n-propylamine	1.068	0.989	7.4	106	0.00
16		4-Methylphenol	1.401	1.324	5.5	111	0.00
17		Hexachloroethane	0.551	0.550	0.2	114	0.00
18	I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
19	S	Nitrobenzene-d5	0.145	0.148	-2.1	110	0.00
20		Nitrobenzene	0.374	0.383	-2.4	112	0.00
21		Isophorone	0.700	0.668	4.6	105	0.00
22	S	2-Nitrophenol-d4	0.155	0.167	-7.7	117	0.00
23	C	2-Nitrophenol	0.170	0.181	-6.5	114	0.00
24		2,4-Dimethylphenol	0.389	0.395	-1.5	110	0.00
25		Bis(2-Chloroethoxy)methane	0.420	0.412	1.9	108	0.00
26	S	2,4-Dichlorophenol-d3	0.325	0.333	-2.5	113	0.00
27	C	2,4-Dichlorophenol	0.324	0.333	-2.8	112	0.00
28		Naphthalene	0.991	0.995	-0.4	110	0.00
29	S	4-Chloroaniline-d4	0.387	0.394	-1.8	107	0.00
30		4-Chloroaniline	0.397	0.401	-1.0	105	0.00
31	C	Hexachlorobutadiene	0.239	0.246	-2.9	113	0.00
32		Caprolactam	0.110	0.095	13.6	99	0.00
33	C	4-Chloro-3-methylphenol	0.354	0.345	2.5	107	0.00
34		2-Methylnaphthalene	0.764	0.751	1.7	108	0.00
35	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.00
36		1,2,4,5-Tetrachlorobenzene	0.707	0.787	-11.3	109	0.00
37		Hexachlorocyclopentadiene	0.450	0.363	19.3	80	0.00
38	C	2,4,6-Trichlorophenol	0.454	0.487	-7.3	106	0.00
39		2,4,5-Trichlorophenol	0.475	0.506	-6.5	104	0.00
40		1,1'-Biphenyl	1.605	1.711	-6.6	106	0.00
41		2-Chloronaphthalene	1.259	1.326	-5.3	105	0.00
42		2-Nitroaniline	0.361	0.378	-4.7	104	0.00
43	S	Dimethylphthalate-d6	1.653	1.650	0.2	100	0.00
44		Dimethylphthalate	1.652	1.635	1.0	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM081816\
Data File : BM007176.D
Acq On : 17 Aug 2016 12:39
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02056

Quant Time: Aug 17 13:54:30 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM081116.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 17 04:01:43 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	0.321	0.337	-5.0	105	0.00
46	S	Acenaphthylene-d8	1.985	2.054	-3.5	102	0.00
47		Acenaphthylene	2.043	2.120	-3.8	103	0.00
48		3-Nitroaniline	0.317	0.335	-5.7	100	0.00
49	C	Acenaphthene	1.327	1.354	-2.0	102	0.00
50		2,4-Dinitrophenol	0.190	0.142	25.3#	88	0.00
51	S	4-Nitrophenol-d4	0.296	0.278	6.1	95	0.00
52		4-Nitrophenol	0.301	0.287	4.7	97	0.00
53		Dibenzofuran	1.969	2.014	-2.3	103	0.00
54		2,4-Dinitrotoluene	0.474	0.487	-2.7	102	0.00
55		2,3,4,6-Tetrachlorophenol	0.459	0.472	-2.8	102	0.00
56		Diethylphthalate	1.742	1.636	6.1	97	0.00
57	S	Fluorene-d10	1.445	1.450	-0.3	101	0.00
58		Fluorene	1.588	1.590	-0.1	99	0.00
59		4-Chlorophenyl-phenylether	0.838	0.832	0.7	99	0.00
60		4-Nitroaniline	0.372	0.360	3.2	95	0.00
61	I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
62	S	4,6-Dinitro-2-methylphenol-	0.111	0.095	14.4	90	0.00
63		4,6-Dinitro-2-methylphenol	0.120	0.104	13.3	89	0.00
64		N-Nitrosodiphenylamine	0.566	0.580	-2.5	97	0.00
65		4-Bromophenyl-phenylether	0.227	0.234	-3.1	101	0.00
66		Hexachlorobenzene	0.256	0.261	-2.0	99	0.00
67		Atrazine	0.228	0.227	0.4	94	0.00
68	C	Pentachlorophenol	0.158	0.154	2.5	98	0.00
69		Phenanthrene	1.076	1.087	-1.0	98	0.00
70	S	Anthracene-d10	0.927	0.943	-1.7	98	0.00
71		Anthracene	1.104	1.113	-0.8	96	0.00
72		Carbazole	0.951	0.992	-4.3	96	0.00
73		Di-n-butylphthalate	1.163	1.160	0.3	95	0.00
74	C	Fluoranthene	1.305	1.406	-7.7	98	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	95	0.00
76	S	Pyrene-d10	0.844	0.871	-3.2	97	0.00
77		Pyrene	1.106	1.121	-1.4	95	0.00
78		Butylbenzylphthalate	0.430	0.432	-0.5	93	0.00
79		3,3'-Dichlorobenzidine	0.386	0.391	-1.3	82	0.00
80		Benzo(a)anthracene	1.182	1.207	-2.1	95	0.00
81		Bis(2-ethylhexyl)phthalate	0.636	0.603	5.2	87	0.00
82		Chrysene	1.080	1.081	-0.1	93	0.00
83	I	Perylene-d12	1.000	1.000	0.0	93	0.00
84		Di-n-octyl phthalate	1.032	1.038	-0.6	89	0.00
85		Benzo(b)fluoranthene	1.205	1.262	-4.7	95	0.00
86		Benzo(k)fluoranthene	1.131	1.125	0.5	91	0.00
87	S	Benzo(a)pyrene-d12	0.915	0.948	-3.6	95	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM081816\
Data File : BM007176.D
Acq On : 17 Aug 2016 12:39
Operator : UM/SJ
Sample : SSTDCCC020
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTD02056

Quant Time: Aug 17 13:54:30 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM081116.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Aug 17 04:01:43 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	1.142	1.172	-2.6	93	0.00
89	Indeno(1,2,3-cd)pyrene	1.400	1.419	-1.4	92	0.00
90	Dibenzo(a,h)anthracene	1.175	1.187	-1.0	92	0.00
91	Benzo(g,h,i)perylene	1.183	1.177	0.5	92	0.00

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