Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF110918\
Data File : BF110615.D
Acq On : 9 Nov 2018 9:29
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 09 15:48:17 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 09 15:13:13 2018
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	105	0.00
2		1,4-Dioxane	0.613	0.617	-0.7	104	0.00
3		Pyridine	1.324	1.405	-6.1	103	0.00
4		n-Nitrosodimethylamine	0.568	0.590	-3.9	103	0.00
5	S	2-Fluorophenol	1.231	1.256	-2.0	103	0.00
6		Aniline	2.053	2.033	1.0	104	0.00
7	S	Phenol-d6	1.575	1.576	-0.1	104	0.00
8		2-Chlorophenol	1.443	1.437	0.4	103	0.00
9		Benzaldehyde	0.893	0.842	5.7	101	0.00
10	C	Phenol	1.826	1.814	0.7	102	0.00
11		bis(2-Chloroethyl)ether	1.368	1.349	1.4	103	0.00
12		1,3-Dichlorobenzene	1.579	1.568	0.7	102	0.00
13	C	1,4-Dichlorobenzene	1.604	1.567	2.3	102	0.00
14		1,2-Dichlorobenzene	1.492	1.480	0.8	104	0.00
15		Benzyl Alcohol	1.232	1.238	-0.5	103	0.00
16		2,2'-oxybis(1-Chloropropane	2.149	2.121	1.3	102	0.00
17		2-Methylphenol	1.149	1.134	1.3	103	0.00
18		Hexachloroethane	0.592	0.579	2.2	101	0.00
19	P	n-Nitroso-di-n-propylamine	1.005	0.994	1.1	104	0.00
20		3+4-Methylphenols	1.475	1.463	0.8	104	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
22		Acetophenone	0.466	0.467	-0.2	102	0.00
23	S	Nitrobenzene-d5	0.347	0.353	-1.7	103	0.00
24		Nitrobenzene	0.356	0.358	-0.6	102	0.00
25		Isophorone	0.569	0.568	0.2	103	0.00
26	C	2-Nitrophenol	0.178	0.185	-3.9	103	0.00
27		2,4-Dimethylphenol	0.270	0.274	-1.5	103	0.00
28		bis(2-Chloroethoxy)methane	0.385	0.382	0.8	101	0.00
29	C	2,4-Dichlorophenol	0.278	0.285	-2.5	102	0.00
30		1,2,4-Trichlorobenzene	0.314	0.315	-0.3	101	0.00
31		Naphthalene	0.939	0.932	0.7	102	0.00
32		Benzoic acid	0.191	0.208	-8.9	103	0.00
33		4-Chloroaniline	0.425	0.415	2.4	103	0.00
34	C	Hexachlorobutadiene	0.179	0.176	1.7	101	0.00
35		Caprolactam	0.093	0.095	-2.2	103	0.00
36	C	4-Chloro-3-methylphenol	0.300	0.304	-1.3	102	0.00
37		2-Methylnaphthalene	0.615	0.613	0.3	102	0.00
38		1-Methylnaphthalene	0.592	0.591	0.2	102	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	102	0.00
40		1,2,4,5-Tetrachlorobenzene	0.633	0.634	-0.2	101	0.00
41	P	Hexachlorocyclopentadiene	0.203	0.195	3.9	92	0.00
42	S	2,4,6-Tribromophenol	0.202	0.195	3.5	98	0.00
43	C	2,4,6-Trichlorophenol	0.398	0.398	0.0	99	0.00
44		2,4,5-Trichlorophenol	0.424	0.425	-0.2	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF110918\
Data File : BF110615.D
Acq On : 9 Nov 2018 9:29
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 09 15:48:17 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 09 15:13:13 2018
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.400	1.360	2.9	100	0.00
46		1,1'-Biphenyl	1.866	1.841	1.3	101	0.00
47		2-Chloronaphthalene	1.344	1.332	0.9	102	0.00
48		2-Nitroaniline	0.414	0.415	-0.2	100	0.00
49		Acenaphthylene	1.936	1.907	1.5	100	0.00
50		Dimethylphthalate	1.492	1.427	4.4	98	0.00
51		2,6-Dinitrotoluene	0.328	0.322	1.8	99	0.00
52	C	Acenaphthene	1.223	1.190	2.7	100	0.00
53		3-Nitroaniline	0.380	0.382	-0.5	102	0.00
54	P	2,4-Dinitrophenol	0.116	0.113	2.6	93	0.00
55		Dibenzofuran	1.790	1.736	3.0	99	0.00
56	P	4-Nitrophenol	0.252	0.252	0.0	97	0.00
57		2,4-Dinitrotoluene	0.427	0.427	0.0	100	0.00
58		Fluorene	1.375	1.329	3.3	100	0.00
59		2,3,4,6-Tetrachlorophenol	0.334	0.334	0.0	101	0.00
60		Diethylphthalate	1.476	1.409	4.5	99	0.00
61		4-Chlorophenyl-phenylether	0.712	0.691	2.9	99	0.00
62		4-Nitroaniline	0.383	0.376	1.8	99	0.00
63		Azobenzene	1.440	1.404	2.5	99	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
65		4,6-Dinitro-2-methylphenol	0.101	0.097	4.0	92	0.00
66	c	n-Nitrosodiphenylamine	0.674	0.678	-0.6	100	0.00
67		4-Bromophenyl-phenylether	0.221	0.217	1.8	98	0.00
68		Hexachlorobenzene	0.233	0.233	0.0	100	0.00
69		Atrazine	0.206	0.193	6.3	93	0.00
70	C	Pentachlorophenol	0.124	0.127	-2.4	97	0.00
71		Phenanthrene	1.071	1.059	1.1	100	0.00
72		Anthracene	1.081	1.069	1.1	99	0.00
73		Carbazole	1.038	1.043	-0.5	100	0.00
74		Di-n-butylphthalate	1.217	1.187	2.5	96	0.00
75	C	Fluoranthene	1.074	1.066	0.7	100	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
77		Benzidine	0.550	0.509	7.5	98	0.00
78		Pyrene	1.540	1.495	2.9	100	0.00
79	S	Terphenyl-d14	1.068	1.012	5.2	100	0.00
80		Butylbenzylphthalate	0.663	0.649	2.1	98	0.00
81		Benzo(a)anthracene	1.195	1.185	0.8	102	0.00
82		3,3'-Dichlorobenzidine	0.400	0.392	2.0	102	0.00
83		Chrysene	1.139	1.126	1.1	105	0.00
84		Bis(2-ethylhexyl)phthalate	0.822	0.807	1.8	101	0.00
85	c	Di-n-octyl phthalate	1.209	1.219	-0.8	107	0.00
86		Indeno(1,2,3-cd)pyrene	0.817	0.744	8.9	96	0.00
87	I	Perylene-d12	1.000	1.000	0.0	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF110918\
Data File : BF110615.D
Acq On : 9 Nov 2018 9:29
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 09 15:48:17 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 09 15:13:13 2018
Response via : Initial Calibration

Min. RRF : 0.000 Min. Rel. Area : 50% Max. R.T. Dev 0.50min
Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.196	1.231	-2.9	98	0.00
89	Benzo(k)fluoranthene	1.182	1.256	-6.3	108	0.00
90 C	Benzo(a)pyrene	1.087	1.095	-0.7	99	0.00
91	Dibenzo(a,h)anthracene	0.920	0.914	0.7	95	-0.01
92	Benzo(g,h,i)perylene	0.913	0.914	-0.1	95	0.00

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