Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112018\
Data File : BF110911.D
Acq On : 20 Nov 2018 23:52
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 21 05:58:59 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 13 12:43:10 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	65	0.00
2		1,4-Dioxane	0.613	0.604	1.5	63	-0.06
3		Pyridine	1.324	1.333	-0.7	61	-0.04
4		n-Nitrosodimethylamine	0.568	0.587	-3.3	63	-0.04
5	S	2-Fluorophenol	1.231	1.301	-5.7	66	0.00
6		Aniline	2.053	1.989	3.1	63	0.00
7	S	Phenol-d6	1.575	1.618	-2.7	66	0.00
8		2-Chlorophenol	1.443	1.481	-2.6	66	0.00
9		Benzaldehyde	0.893	0.904	-1.2	68	0.00
10	C	Phenol	1.826	1.852	-1.4	65	0.00
11		bis(2-Chloroethyl)ether	1.368	1.349	1.4	64	0.00
12		1,3-Dichlorobenzene	1.579	1.598	-1.2	65	0.00
13	C	1,4-Dichlorobenzene	1.604	1.633	-1.8	66	0.00
14		1,2-Dichlorobenzene	1.492	1.522	-2.0	66	-0.01
15		Benzyl Alcohol	1.232	1.252	-1.6	64	0.00
16		2,2'-oxybis(1-Chloropropane	2.149	2.150	-0.0	64	0.00
17		2-Methylphenol	1.149	1.139	0.9	64	0.00
18		Hexachloroethane	0.592	0.510	13.9	55	-0.01
19	P	n-Nitroso-di-n-propylamine	1.005	1.024	-1.9	66	0.00
20		3+4-Methylphenols	1.475	1.498	-1.6	66	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	62	0.00
22		Acetophenone	0.466	0.499	-7.1	66	0.00
23	S	Nitrobenzene-d5	0.347	0.362	-4.3	64	0.00
24		Nitrobenzene	0.356	0.370	-3.9	64	0.00
25		Isophorone	0.569	0.572	-0.5	63	0.00
26	C	2-Nitrophenol	0.178	0.177	0.6	60	0.00
27		2,4-Dimethylphenol	0.270	0.279	-3.3	64	0.00
28		bis(2-Chloroethoxy)methane	0.385	0.401	-4.2	64	0.00
29	C	2,4-Dichlorophenol	0.278	0.288	-3.6	63	0.00
30		1,2,4-Trichlorobenzene	0.314	0.320	-1.9	62	0.00
31		Naphthalene	0.939	0.949	-1.1	63	0.00
32		Benzoic acid	0.191	0.116	39.3#	35#	-0.02
33		4-Chloroaniline	0.425	0.412	3.1	62	0.00
34	C	Hexachlorobutadiene	0.179	0.184	-2.8	64	0.00
35		Caprolactam	0.093	0.087	6.5	58	-0.01
36	C	4-Chloro-3-methylphenol	0.300	0.297	1.0	60	0.00
37		2-Methylnaphthalene	0.615	0.597	2.9	60	0.00
38		1-Methylnaphthalene	0.592	0.571	3.5	60	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	54	0.00
40		1,2,4,5-Tetrachlorobenzene	0.633	0.692	-9.3	59	0.00
41	P	Hexachlorocyclopentadiene	0.203	0.000#	100.0#	0#	-9.09#
42	S	2,4,6-Tribromophenol	0.202	0.176	12.9	47#	0.00
43	C	2,4,6-Trichlorophenol	0.398	0.425	-6.8	56	0.00
44		2,4,5-Trichlorophenol	0.424	0.439	-3.5	55	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112018\
Data File : BF110911.D
Acq On : 20 Nov 2018 23:52
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 21 05:58:59 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 13 12:43:10 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	1.562	-11.6	61	0.00
46		1,1'-Biphenyl	1.866	2.040	-9.3	59	0.00
47		2-Chloronaphthalene	1.344	1.420	-5.7	58	0.00
48		2-Nitroaniline	0.414	0.453	-9.4	58	0.00
49		Acenaphthylene	1.936	1.947	-0.6	54	-0.01
50		Dimethylphthalate	1.492	1.607	-7.7	59	0.00
51		2,6-Dinitrotoluene	0.328	0.325	0.9	53	0.00
52	C	Acenaphthene	1.223	1.232	-0.7	55	-0.01
53		3-Nitroaniline	0.380	0.397	-4.5	57	0.00
54	P	2,4-Dinitrophenol	0.116	0.003#	97.4#	1#	0.03
55		Dibenzofuran	1.790	1.734	3.1	53	0.00
56	P	4-Nitrophenol	0.252	0.180	28.6#	37#	0.01
57		2,4-Dinitrotoluene	0.427	0.370	13.3	46#	0.00
58		Fluorene	1.375	1.288	6.3	51	0.00
59		2,3,4,6-Tetrachlorophenol	0.334	0.293	12.3	47#	0.00
60		Diethylphthalate	1.476	1.560	-5.7	58	0.00
61		4-Chlorophenyl-phenylether	0.712	0.717	-0.7	55	0.00
62		4-Nitroaniline	0.383	0.337	12.0	47#	0.00
63		Azobenzene	1.440	1.354	6.0	51	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	47#	0.00
65		4,6-Dinitro-2-methylphenol	0.101	0.011	89.1#	5#	0.01
66	c	n-Nitrosodiphenylamine	0.674	0.747	-10.8	52	-0.01
67		4-Bromophenyl-phenylether	0.221	0.237	-7.2	51	0.00
68		Hexachlorobenzene	0.233	0.237	-1.7	48#	0.00
69		Atrazine	0.206	0.205	0.5	47#	0.00
70	C	Pentachlorophenol	0.124	0.102	17.7	37#	0.00
71		Phenanthrene	1.071	1.080	-0.8	49#	0.00
72		Anthracene	1.081	1.090	-0.8	48#	0.00
73		Carbazole	1.038	1.063	-2.4	48#	0.00
74		Di-n-butylphthalate	1.217	1.287	-5.8	49#	0.00
75	C	Fluoranthene	1.074	1.130	-5.2	50	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	54	0.00
77		Benzidine	0.550	0.715	-30.0#	72	0.00
78		Pyrene	1.540	1.458	5.3	51	0.00
79	S	Terphenyl-d14	1.068	1.038	2.8	54	0.00
80		Butylbenzylphthalate	0.663	0.683	-3.0	54	0.00
81		Benzo(a)anthracene	1.195	1.194	0.1	54	0.00
82		3,3'-Dichlorobenzidine	0.400	0.467	-16.8	64	0.00
83		Chrysene	1.139	1.160	-1.8	57	0.00
84		Bis(2-ethylhexyl)phthalate	0.822	0.944	-14.8	62	0.00
85	c	Di-n-octyl phthalate	1.209	1.563	-29.3#	72	0.00
86		Indeno(1,2,3-cd)pyrene	0.817	0.733	10.3	49#	0.02
87	I	Perylene-d12	1.000	1.000	0.0	51	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF112018\
Data File : BF110911.D
Acq On : 20 Nov 2018 23:52
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 21 05:58:59 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF110818.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 13 12:43:10 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	1.188	0.7	50#	0.00
89	Benzo(k)fluoranthene	1.182	1.317	-11.4	59	0.00
90 C	Benzo(a)pyrene	1.087	1.094	-0.6	52	0.00
91	Dibenzo(a,h)anthracene	0.920	0.923	-0.3	50	0.01
92	Benzo(g,h,i)perylene	0.913	0.892	2.3	49#	0.03

(#) = Out of Range SPCC's out = 2 CCC's out = 1