Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
Data File : BF085010.D
Acq On : 11 Feb 2016 11:09
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 11 16:13:02 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 10 13:28:51 2016
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	114	-0.01
2		1,4-Dioxane	0.562	0.516	8.2	107	-0.03
3		Pyridine	1.465	1.448	1.2	119	-0.03
4		n-Nitrosodimethylamine	0.758	0.758	0.0	112	-0.03
5	S	2-Fluorophenol	1.284	1.176	8.4	111	-0.01
6		Aniline	1.948	1.967	-1.0	117	-0.02
7	S	Phenol-d6	1.592	1.475	7.3	113	-0.01
8		2-Chlorophenol	1.453	1.359	6.5	105	-0.01
9		Benzaldehyde	0.940	0.836	11.1	99	-0.01
10	C	Phenol	1.783	1.588	10.9	116	-0.01
11		bis(2-Chloroethyl)ether	1.391	1.424	-2.4	127	-0.01
12		1,3-Dichlorobenzene	1.536	1.386	9.8	109	-0.01
13	C	1,4-Dichlorobenzene	1.535	1.405	8.5	110	-0.01
14		1,2-Dichlorobenzene	1.430	1.419	0.8	110	-0.01
15		Benzyl Alcohol	1.099	1.134	-3.2	121	-0.01
16		2,2'-oxybis(1-Chloropropane	2.339	2.237	4.4	116	-0.01
17		2-Methylphenol	1.149	1.129	1.7	107	-0.01
18		Hexachloroethane	0.591	0.573	3.0	109	-0.01
19	P	n-Nitroso-di-n-propylamine	0.998	0.921	7.7	113	-0.01
20		3+4-Methylphenols	1.427	1.484	-4.0	121	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	116	-0.01
22		Acetophenone	0.527	0.479	9.1	108	-0.01
23	S	Nitrobenzene-d5	0.355	0.348	2.0	113	-0.01
24		Nitrobenzene	0.380	0.372	2.1	127	-0.01
25		Isophorone	0.690	0.659	4.5	110	-0.01
26	C	2-Nitrophenol	0.188	0.198	-5.3	125	-0.01
27		2,4-Dimethylphenol	0.326	0.299	8.3	114	-0.01
28		bis(2-Chloroethoxy)methane	0.410	0.358	12.7	106	-0.01
29	C	2,4-Dichlorophenol	0.279	0.255	8.6	103	-0.01
30		1,2,4-Trichlorobenzene	0.315	0.306	2.9	116	-0.01
31		Naphthalene	1.003	0.957	4.6	119	-0.01
32		Benzoic acid	0.209	0.237	-13.4	132	0.00
33		4-Chloroaniline	0.415	0.428	-3.1	132	-0.01
34	C	Hexachlorobutadiene	0.182	0.171	6.0	108	-0.01
35		Caprolactam	0.086	0.090	-4.7	108	0.00
36	C	4-Chloro-3-methylphenol	0.318	0.291	8.5	115	-0.01
37		2-Methylnaphthalene	0.628	0.617	1.8	109	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	110	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.635	0.619	2.5	118	-0.01
40	P	Hexachlorocyclopentadiene	0.223	0.234	-4.9	115	-0.01
41	S	2,4,6-Tribromophenol	0.209	0.212	-1.4	107	-0.01
42	C	2,4,6-Trichlorophenol	0.409	0.402	1.7	107	-0.01
43		2,4,5-Trichlorophenol	0.416	0.418	-0.5	115	-0.01
44	S	2-Fluorobiphenyl	1.321	1.107	16.2	104	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
Data File : BF085010.D
Acq On : 11 Feb 2016 11:09
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 11 16:13:02 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 10 13:28:51 2016
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		1,1'-Biphenyl	1.786	1.847	-3.4	114	-0.01
46		2-Chloronaphthalene	1.316	1.300	1.2	112	-0.01
47		2-Nitroaniline	0.417	0.464	-11.3	139	-0.01
48		Acenaphthylene	1.996	1.952	2.2	108	-0.01
49		Dimethylphthalate	1.523	1.418	6.9	102	-0.01
50		2,6-Dinitrotoluene	0.333	0.327	1.8	103	-0.01
51	C	Acenaphthene	1.299	1.289	0.8	110	-0.01
52		3-Nitroaniline	0.374	0.327	12.6	100	-0.01
53	P	2,4-Dinitrophenol	0.133	0.147	-10.5	114	-0.01
54		Dibenzofuran	1.587	1.574	0.8	109	-0.01
55	P	4-Nitrophenol	0.275	0.238	13.5	87	-0.01
56		2,4-Dinitrotoluene	0.424	0.441	-4.0	113	-0.01
57		Fluorene	1.326	1.376	-3.8	112	-0.01
58		2,3,4,6-Tetrachlorophenol	0.317	0.333	-5.0	130	-0.01
59		Diethylphthalate	1.487	1.312	11.8	101	-0.01
60		4-Chlorophenyl-phenylether	0.621	0.526	15.3	100	-0.01
61		4-Nitroaniline	0.364	0.340	6.6	106	-0.01
62		Azobenzene	1.430	1.347	5.8	106	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	114	-0.01
64		4,6-Dinitro-2-methylphenol	0.123	0.138	-12.2	121	0.00
65	c	n-Nitrosodiphenylamine	0.635	0.575	9.4	96	-0.01
66		4-Bromophenyl-phenylether	0.221	0.222	-0.5	107	-0.01
67		Hexachlorobenzene	0.241	0.217	10.0	108	-0.01
68		Atrazine	0.201	0.190	5.5	115	-0.01
69	C	Pentachlorophenol	0.113	0.119	-5.3	115	-0.01
70		Phenanthrene	1.109	0.964	13.1	106	-0.01
71		Anthracene	1.118	1.145	-2.4	109	-0.01
72		Carbazole	0.975	0.971	0.4	105	-0.01
73		Di-n-butylphthalate	1.241	1.198	3.5	115	0.00
74	C	Fluoranthene	1.123	1.114	0.8	108	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	92	-0.01
76		Benzidine	0.597	0.702	-17.6	90	-0.01
77		Pyrene	1.531	1.654	-8.0	95	-0.01
78	S	Terphenyl-d14	0.883	0.962	-8.9	104	-0.01
79		Butylbenzylphthalate	0.695	0.795	-14.4	98	-0.01
80		Benzo(a)anthracene	1.241	1.278	-3.0	94	-0.01
81		3,3'-Dichlorobenzidine	0.440	0.495	-12.5	94	-0.01
82		Chrysene	1.204	1.266	-5.1	94	0.00
83		Bis(2-ethylhexyl)phthalate	0.864	0.840	2.8	87	-0.01
84	c	Di-n-octyl phthalate	1.426	1.361	4.6	85	0.00
85		Indeno(1,2,3-cd)pyrene	0.947	1.159	-22.4	114	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	101	-0.01
87		Benzo(b)fluoranthene	1.344	1.243	7.5	93	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF021116\
Data File : BF085010.D
Acq On : 11 Feb 2016 11:09
Operator : SJ/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 11 16:13:02 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF020116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 10 13:28:51 2016
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(k)fluoranthene	1.111	0.949	14.6	89	-0.01
89 C	Benzo(a)pyrene	1.145	1.052	8.1	92	-0.01
90	Dibenzo(a,h)anthracene	0.964	1.061	-10.1	113	-0.01
91	Benzo(g,h,i)perylene	0.971	1.131	-16.5	119	0.00

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