Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF061115\
Data File : BF079896.D
Acq On : 11 Jun 2015 15:36
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF061115

Quant Time: Jun 12 09:14:57 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 08:55:39 2015
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	116	-0.14
2		1,4-Dioxane	0.496	0.508	-2.4	116	0.38
3		Pyridine	1.272	1.562	-22.8	152#	0.75#
4		n-Nitrosodimethylamine	0.627	0.596	4.9	120	0.75#
5	S	2-Fluorophenol	1.091	1.110	-1.7	116	0.88#
6		Aniline	2.151	2.284	-6.2	117	0.19
7	S	Phenol-d6	1.486	1.616	-8.7	117	0.26
8		2-Chlorophenol	1.326	1.421	-7.2	118	0.08
9		Benzaldehyde	0.854	0.910	-6.6	131	0.26
10	C	Phenol	1.634	1.779	-8.9	117	0.24
11		bis(2-Chloroethyl)ether	1.315	1.385	-5.3	113	0.14
12		1,3-Dichlorobenzene	1.546	1.586	-2.6	114	-0.08
13	C	1,4-Dichlorobenzene	1.606	1.681	-4.7	118	-0.15
14		1,2-Dichlorobenzene	1.427	1.504	-5.4	119	-0.31
15		Benzyl Alcohol	1.191	1.259	-5.7	116	-0.26
16		2,2'-oxybis(1-Chloropropane	1.948	1.977	-1.5	112	-0.42
17		2-Methylphenol	1.158	1.257	-8.5	120	-0.38
18		Hexachloroethane	0.497	0.521	-4.8	117	-0.67#
19	P	n-Nitroso-di-n-propylamine	1.062	1.089	-2.5	112	-0.56#
20		3+4-Methylphenols	1.480	1.602	-8.2	117	-0.56#
21	I	Naphthalene-d8	1.000	1.000	0.0	114	-1.64#
22		Acetophenone	0.472	0.507	-7.4	118	-0.57#
23	S	Nitrobenzene-d5	0.303	0.341	-12.5	125	-0.74#
24		Nitrobenzene	0.322	0.333	-3.4	114	-0.78#
25		Isophorone	0.705	0.737	-4.5	114	-1.05#
26	C	2-Nitrophenol	0.100	0.116	-16.0	126	-1.15#
27		2,4-Dimethylphenol	0.316	0.320	-1.3	117	-1.22#
28		bis(2-Chloroethoxy)methane	0.419	0.426	-1.7	113	-1.34#
29	C	2,4-Dichlorophenol	0.265	0.275	-3.8	118	-1.47#
30		1,2,4-Trichlorobenzene	0.337	0.345	-2.4	117	-1.56#
31		Naphthalene	1.093	1.123	-2.7	115	-1.67#
32		Benzoic acid	0.107	0.114	-6.5	113	-1.35#
33		4-Chloroaniline	0.424	0.431	-1.7	111	-1.76#
34	C	Hexachlorobutadiene	0.187	0.187	0.0	117	-1.82#
35		Caprolactam	0.084	0.094	-11.9	111	-2.19#
36	C	4-Chloro-3-methylphenol	0.302	0.324	-7.3	114	-2.45#
37		2-Methylnaphthalene	0.739	0.769	-4.1	117	-2.60#
38	I	Acenaphthene-d10	1.000	1.000	0.0	115	-3.75#
39		1,2,4,5-Tetrachlorobenzene	0.670	0.664	0.9	116	-2.80#
40	P	Hexachlorocyclopentadiene	0.211	0.236	-11.8	117	-2.79#
41	S	2,4,6-Tribromophenol	0.182	0.200	-9.9	111	-4.45#
42	C	2,4,6-Trichlorophenol	0.413	0.441	-6.8	117	-2.96#
43		2,4,5-Trichlorophenol	0.456	0.458	-0.4	110	-3.01#
44	S	2-Fluorobiphenyl	1.428	1.355	5.1	114	-3.06#

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF061115\
Data File : BF079896.D
Acq On : 11 Jun 2015 15:36
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF061115

Quant Time: Jun 12 09:14:57 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 08:55:39 2015
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	1.619	1.586	2.0	114	-3.16#
46		2-Chloronaphthalene	1.378	1.339	2.8	114	-3.17#
47		2-Nitroaniline	0.256	0.286	-11.7	119	-3.32#
48		Acenaphthylene	2.326	2.325	0.0	114	-3.60#
49		Dimethylphthalate	1.602	1.532	4.4	111	-3.53#
50		2,6-Dinitrotoluene	0.250	0.295	-18.0	117	-3.58#
51	C	Acenaphthene	1.322	1.296	2.0	114	-3.78#
52		3-Nitroaniline	0.319	0.352	-10.3	116	-3.75#
53	P	2,4-Dinitrophenol	0.037	0.041#	-10.8	121	-3.86#
54		Dibenzofuran	1.932	1.931	0.1	114	-3.95#
55	P	4-Nitrophenol	0.215	0.262	-21.9	123	-3.97#
56		2,4-Dinitrotoluene	0.314	0.391	-24.5	120	-3.96#
57		Fluorene	1.684	1.725	-2.4	113	-4.25#
58		2,3,4,6-Tetrachlorophenol	0.345	0.369	-7.0	113	-4.07#
59		Diethylphthalate	1.602	1.605	-0.2	111	-4.19#
60		4-Chlorophenyl-phenylether	0.752	0.735	2.3	113	-4.27#
61		4-Nitroaniline	0.326	0.374	-14.7	113	-4.30#
62		Azobenzene	1.542	1.521	1.4	111	-4.40#
63	I	Phenanthrene-d10	1.000	1.000	0.0	111	-4.99#
64		4,6-Dinitro-2-methylphenol	0.038	0.047	-23.7	130	-4.32#
65	c	n-Nitrosodiphenylamine	0.646	0.671	-3.9	112	-4.37#
66		4-Bromophenyl-phenylether	0.224	0.244	-8.9	116	-4.66#
67		Hexachlorobenzene	0.244	0.262	-7.4	114	-4.69#
68		Atrazine	0.199	0.213	-7.0	109	-4.80#
69	C	Pentachlorophenol	0.095	0.098	-3.2	110	-4.87#
70		Phenanthrene	1.192	1.215	-1.9	109	-5.00#
71		Anthracene	1.153	1.194	-3.6	111	-5.05#
72		Carbazole	1.119	1.120	-0.1	104	-5.18#
73		Di-n-butylphthalate	1.219	1.286	-5.5	111	-5.41#
74	C	Fluoranthene	1.318	1.385	-5.1	110	-5.76#
75	I	Chrysene-d12	1.000	1.000	0.0	109	-6.19#
76		Benzidine	0.554	0.608	-9.7	123	-5.85#
77		Pyrene	1.256	1.255	0.1	112	-5.86#
78	S	Terphenyl-d14	0.867	0.859	0.9	111	-5.92#
79		Butylbenzylphthalate	0.424	0.461	-8.7	111	-6.07#
80		Benzo(a)anthracene	1.270	1.209	4.8	109	-6.18#
81		3,3'-Dichlorobenzidine	0.406	0.414	-2.0	108	-6.18#
82		Chrysene	1.229	1.136	7.6	108	-6.19#
83		Bis(2-ethylhexyl)phthalate	0.679	0.731	-7.7	110	-6.14#
84	c	Di-n-octyl phthalate	1.107	1.197	-8.1	108	-6.18#
85		Indeno(1,2,3-cd)pyrene	1.330	1.338	-0.6	112	-6.75#
86	I	Perylene-d12	1.000	1.000	0.0	109	-6.45#
87		Benzo(b)fluoranthene	1.225	1.198	2.2	110	-6.35#

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF061115\
Data File : BF079896.D
Acq On : 11 Jun 2015 15:36
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF061115

Quant Time: Jun 12 09:14:57 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 12 08:55:39 2015
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	1.250	1.317	-5.4	109	-6.35#
89 C	Benzo(a)pyrene	1.165	1.181	-1.4	110	-6.44#
90	Dibenzo(a,h)anthracene	1.146	1.139	0.6	112	-6.73#
91	Benzo(g,h,i)perylene	1.185	1.179	0.5	112	-6.86#

(#) = Out of Range SPCC's out = 1 CCC's out = 0