Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF112215\
Data File : BF083151.D
Acq On : 22 Nov 2015 11:23
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 23 03:57:43 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Nov 22 09:22:46 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	110	-0.01
2		1,4-Dioxane	0.655	0.648	1.1	114	-0.06
3		Pyridine	1.731	1.843	-6.5	119	-0.27
4		n-Nitrosodimethylamine	0.798	0.894	-12.0	113	-0.08
5	S	2-Fluorophenol	1.323	1.313	0.8	112	-0.03
6		Aniline	2.092	2.280	-9.0	116	-0.02
7	S	Phenol-d6	1.712	1.697	0.9	115	-0.02
8		2-Chlorophenol	1.429	1.449	-1.4	115	-0.01
9		Benzaldehyde	0.894	0.906	-1.3	115	-0.02
10	C	Phenol	2.071	2.059	0.6	123	-0.02
11		bis(2-Chloroethyl)ether	1.479	1.608	-8.7	123	-0.01
12		1,3-Dichlorobenzene	1.632	1.633	-0.1	114	0.00
13	C	1,4-Dichlorobenzene	1.651	1.618	2.0	112	0.00
14		1,2-Dichlorobenzene	1.500	1.386	7.6	102	-0.01
15		Benzyl Alcohol	1.430	1.328	7.1	101	-0.02
16		2,2'-oxybis(1-Chloropropane	2.306	2.571	-11.5	127	-0.01
17		2-Methylphenol	1.183	1.160	1.9	112	-0.02
18		Hexachloroethane	0.581	0.584	-0.5	111	-0.01
19	P	n-Nitroso-di-n-propylamine	1.191	1.160	2.6	114	-0.02
20		3+4-Methylphenols	1.495	1.538	-2.9	114	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	115	-0.01
22		Acetophenone	0.574	0.545	5.1	115	-0.02
23	S	Nitrobenzene-d5	0.438	0.422	3.7	113	-0.01
24		Nitrobenzene	0.468	0.462	1.3	116	-0.01
25		Isophorone	0.829	0.818	1.3	116	-0.03
26	C	2-Nitrophenol	0.206	0.205	0.5	107	-0.01
27		2,4-Dimethylphenol	0.330	0.335	-1.5	123	-0.01
28		bis(2-Chloroethoxy)methane	0.482	0.456	5.4	108	-0.02
29	C	2,4-Dichlorophenol	0.313	0.309	1.3	113	-0.01
30		1,2,4-Trichlorobenzene	0.344	0.325	5.5	111	-0.01
31		Naphthalene	1.079	1.063	1.5	116	-0.01
32		Benzoic acid	0.256	0.276	-7.8	115	-0.05
33		4-Chloroaniline	0.419	0.428	-2.1	120	-0.01
34	C	Hexachlorobutadiene	0.203	0.194	4.4	112	0.00
35		Caprolactam	0.100	0.099	1.0	116	-0.07
36	C	4-Chloro-3-methylphenol	0.363	0.355	2.2	116	-0.02
37		2-Methylnaphthalene	0.701	0.636	9.3	109	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	116	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.648	0.620	4.3	115	-0.01
40	P	Hexachlorocyclopentadiene	0.369	0.295	20.1	91	-0.01
41	S	2,4,6-Tribromophenol	0.205	0.189	7.8	105	-0.01
42	C	2,4,6-Trichlorophenol	0.464	0.435	6.3	110	-0.02
43		2,4,5-Trichlorophenol	0.475	0.456	4.0	115	-0.02
44	S	2-Fluorobiphenyl	1.434	1.316	8.2	122	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF112215\
Data File : BF083151.D
Acq On : 22 Nov 2015 11:23
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 23 03:57:43 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Nov 22 09:22:46 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	1.702	1.557	8.5	103	-0.01
46		2-Chloronaphthalene	1.337	1.270	5.0	114	-0.01
47		2-Nitroaniline	0.473	0.476	-0.6	123	-0.01
48		Acenaphthylene	2.072	2.034	1.8	116	-0.01
49		Dimethylphthalate	1.540	1.470	4.5	113	-0.01
50		2,6-Dinitrotoluene	0.345	0.313	9.3	116	-0.01
51	C	Acenaphthene	1.261	1.193	5.4	112	-0.01
52		3-Nitroaniline	0.368	0.345	6.3	106	-0.02
53	P	2,4-Dinitrophenol	0.199	0.163	18.1	94	-0.01
54		Dibenzofuran	1.783	1.696	4.9	111	-0.01
55	P	4-Nitrophenol	0.282	0.255	9.6	108	-0.02
56		2,4-Dinitrotoluene	0.438	0.392	10.5	112	-0.01
57		Fluorene	1.474	1.353	8.2	112	0.00
58		2,3,4,6-Tetrachlorophenol	0.389	0.370	4.9	109	-0.01
59		Diethylphthalate	1.533	1.376	10.2	105	-0.02
60		4-Chlorophenyl-phenylether	0.675	0.590	12.6	108	-0.01
61		4-Nitroaniline	0.363	0.365	-0.6	111	-0.02
62		Azobenzene	1.749	1.645	5.9	114	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	110	-0.01
64		4,6-Dinitro-2-methylphenol	0.138	0.141	-2.2	104	-0.01
65	c	n-Nitrosodiphenylamine	0.680	0.648	4.7	111	-0.01
66		4-Bromophenyl-phenylether	0.222	0.223	-0.5	110	-0.01
67		Hexachlorobenzene	0.244	0.222	9.0	103	-0.01
68		Atrazine	0.198	0.209	-5.6	115	-0.02
69	C	Pentachlorophenol	0.159	0.150	5.7	102	-0.01
70		Phenanthrene	1.152	1.069	7.2	104	-0.01
71		Anthracene	1.176	1.161	1.3	118	0.00
72		Carbazole	1.045	1.071	-2.5	112	-0.01
73		Di-n-butylphthalate	1.275	1.238	2.9	110	-0.01
74	C	Fluoranthene	1.179	1.109	5.9	113	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	106	-0.01
76		Benzidine	0.524	0.590	-12.6	111	-0.01
77		Pyrene	1.366	1.368	-0.1	111	-0.01
78	S	Terphenyl-d14	0.849	0.854	-0.6	108	-0.01
79		Butylbenzylphthalate	0.647	0.683	-5.6	113	-0.01
80		Benzo(a)anthracene	1.237	1.246	-0.7	108	-0.01
81		3,3'-Dichlorobenzidine	0.431	0.409	5.1	103	-0.02
82		Chrysene	1.147	1.105	3.7	109	-0.01
83		Bis(2-ethylhexyl)phthalate	0.845	0.866	-2.5	113	-0.01
84	c	Di-n-octyl phthalate	1.454	1.481	-1.9	112	-0.02
85		Indeno(1,2,3-cd)pyrene	1.043	1.028	1.4	115	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	112	-0.01
87		Benzo(b)fluoranthene	1.367	1.380	-1.0	113	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF112215\
Data File : BF083151.D
Acq On : 22 Nov 2015 11:23
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Nov 23 03:57:43 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Nov 22 09:22:46 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	1.018	0.949	6.8	113	-0.01
89 C	Benzo(a)pyrene	1.128	1.075	4.7	110	-0.02
90	Dibenzo(a,h)anthracene	1.025	0.964	6.0	112	-0.02
91	Benzo(g,h,i)perylene	1.000	0.933	6.7	111	-0.02

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