Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF072315\
Data File : BF080622.D
Acq On : 24 Jul 2015 6:21
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 24 06:47:59 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 24 01:01:51 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	109	0.00
2		1,4-Dioxane	0.552	0.499	9.6	92	0.01
3		Pyridine	1.466	1.473	-0.5	100	0.01
4		n-Nitrosodimethylamine	0.927	0.934	-0.8	102	0.00
5	S	2-Fluorophenol	1.158	1.291	-11.5	113	0.01
6		Aniline	2.069	2.332	-12.7	113	0.00
7	S	Phenol-d6	1.416	1.493	-5.4	102	0.00
8		2-Chlorophenol	1.244	1.435	-15.4	116	0.00
9		Benzaldehyde	1.047	1.097	-4.8	107	0.00
10	C	Phenol	1.694	1.680	0.8	97	0.00
11		bis(2-Chloroethyl)ether	1.176	1.175	0.1	106	0.00
12		1,3-Dichlorobenzene	1.577	1.616	-2.5	108	0.00
13	C	1,4-Dichlorobenzene	1.548	1.650	-6.6	112	0.00
14		1,2-Dichlorobenzene	1.435	1.409	1.8	101	0.00
15		Benzyl Alcohol	1.247	1.470	-17.9	122	0.00
16		2,2'-oxybis(1-Chloropropane	2.384	2.433	-2.1	105	0.00
17		2-Methylphenol	0.980	1.153	-17.7	120	0.01
18		Hexachloroethane	0.535	0.557	-4.1	108	0.00
19	P	n-Nitroso-di-n-propylamine	0.959	1.145	-19.4	126	0.00
20		3+4-Methylphenols	1.355	1.505	-11.1	116	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	111	0.00
22		Acetophenone	0.487	0.543	-11.5	115	0.00
23	S	Nitrobenzene-d5	0.353	0.360	-2.0	104	0.00
24		Nitrobenzene	0.377	0.439	-16.4	118	0.00
25		Isophorone	0.654	0.728	-11.3	115	0.00
26	C	2-Nitrophenol	0.135	0.150	-11.1	120	0.00
27		2,4-Dimethylphenol	0.298	0.337	-13.1	117	0.00
28		bis(2-Chloroethoxy)methane	0.361	0.411	-13.9	124	0.00
29	C	2,4-Dichlorophenol	0.244	0.295	-20.9#	131	0.00
30		1,2,4-Trichlorobenzene	0.317	0.336	-6.0	117	0.00
31		Naphthalene	1.000	1.055	-5.5	114	0.00
32		Benzoic acid	0.127	0.211	-66.1#	184#	0.00
33		4-Chloroaniline	0.392	0.440	-12.2	119	0.00
34	C	Hexachlorobutadiene	0.185	0.193	-4.3	113	0.00
35		Caprolactam	0.066	0.086	-30.3#	127	0.01
36	C	4-Chloro-3-methylphenol	0.290	0.302	-4.1	103	0.00
37		2-Methylnaphthalene	0.560	0.603	-7.7	114	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	117	0.00
39		1,2,4,5-Tetrachlorobenzene	0.676	0.619	8.4	105	0.00
40	P	Hexachlorocyclopentadiene	0.356	0.120	66.3#	38#	0.00
41	S	2,4,6-Tribromophenol	0.158	0.197	-24.7	139	0.00
42	C	2,4,6-Trichlorophenol	0.379	0.401	-5.8	122	0.01
43		2,4,5-Trichlorophenol	0.383	0.427	-11.5	128	0.01
44	S	2-Fluorobiphenyl	1.378	1.264	8.3	108	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF072315\
Data File : BF080622.D
Acq On : 24 Jul 2015 6:21
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 24 06:47:59 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 24 01:01:51 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	1.628	1.484	8.8	106	0.00
46		2-Chloronaphthalene	1.317	1.188	9.8	105	0.00
47		2-Nitroaniline	0.402	0.426	-6.0	123	0.00
48		Acenaphthylene	1.951	1.994	-2.2	120	0.00
49		Dimethylphthalate	1.471	1.524	-3.6	115	0.00
50		2,6-Dinitrotoluene	0.275	0.270	1.8	108	0.00
51	C	Acenaphthene	1.172	1.178	-0.5	118	0.00
52		3-Nitroaniline	0.299	0.354	-18.4	141	0.00
53	P	2,4-Dinitrophenol	0.076	0.027#	64.5#	43#	0.00
54		Dibenzofuran	1.694	1.706	-0.7	115	0.00
55	P	4-Nitrophenol	0.216	0.243	-12.5	134	0.01
56		2,4-Dinitrotoluene	0.345	0.337	2.3	106	0.00
57		Fluorene	1.360	1.380	-1.5	114	0.00
58		2,3,4,6-Tetrachlorophenol	0.298	0.362	-21.5	130	0.00
59		Diethylphthalate	1.373	1.354	1.4	123	0.01
60		4-Chlorophenyl-phenylether	0.611	0.594	2.8	114	0.00
61		4-Nitroaniline	0.301	0.365	-21.3	143	0.00
62		Azobenzene	1.439	1.399	2.8	108	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	123	0.00
64		4,6-Dinitro-2-methylphenol	0.065	0.029	55.4#	58	0.00
65	c	n-Nitrosodiphenylamine	0.656	0.603	8.1	112	0.00
66		4-Bromophenyl-phenylether	0.203	0.185	8.9	111	0.00
67		Hexachlorobenzene	0.230	0.226	1.7	124	0.00
68		Atrazine	0.180	0.188	-4.4	118	0.00
69	C	Pentachlorophenol	0.107	0.152	-42.1#	176#	0.01
70		Phenanthrene	1.129	1.118	1.0	120	0.00
71		Anthracene	1.057	1.146	-8.4	135	0.01
72		Carbazole	0.914	1.090	-19.3	143	0.00
73		Di-n-butylphthalate	1.062	1.297	-22.1	142	0.00
74	C	Fluoranthene	1.010	1.369	-35.5#	162#	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	217#	0.00
76		Benzidine	0.588	0.660	-12.2	235#	0.00
77		Pyrene	1.342	1.038	22.7	168#	0.00
78	S	Terphenyl-d14	0.956	0.750	21.5	170#	0.00
79		Butylbenzylphthalate	0.444	0.497	-11.9	228#	-0.01
80		Benzo(a)anthracene	1.128	1.165	-3.3	220#	0.00
81		3,3'-Dichlorobenzidine	0.322	0.442	-37.3#	274#	0.00
82		Chrysene	1.122	1.148	-2.3	218#	0.00
83		Bis(2-ethylhexyl)phthalate	0.656	0.720	-9.8	229#	-0.01
84	c	Di-n-octyl phthalate	0.829	1.366	-64.8#	329#	-0.01
85		Indeno(1,2,3-cd)pyrene	0.849	1.177	-38.6#	280#	0.02
86	I	Perylene-d12	1.000	1.000	0.0	295#	0.00
87		Benzo(b)fluoranthene	1.216	1.170	3.8	274#	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF072315\
Data File : BF080622.D
Acq On : 24 Jul 2015 6:21
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 24 06:47:59 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 24 01:01:51 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	1.211	1.188	1.9	265#	0.00
89 C	Benzo(a)pyrene	1.079	1.093	-1.3	279#	0.00
90	Dibenzo(a,h)anthracene	1.019	1.031	-1.2	275#	0.01
91	Benzo(g,h,i)perylene	1.041	1.013	2.7	266#	0.01

(#) = Out of Range SPCC's out = 1 CCC's out = 4