Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF020215\
Data File : BF077157.D
Acq On : 2 Feb 2015 14:09
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 03 09:57:22 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 03 00:09:26 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	148	0.00
2		1,4-Dioxane	0.521	0.498	4.4	149	0.00
3		Pyridine	1.348	1.119	17.0	102	0.00
4		n-Nitrosodimethylamine	0.668	0.732	-9.6	161#	0.00
5	S	2-Fluorophenol	1.216	1.168	3.9	135	0.00
6		Aniline	2.124	2.072	2.4	135	0.00
7	S	Phenol-d6	1.535	1.477	3.8	135	0.00
8		2-Chlorophenol	1.402	1.358	3.1	134	0.00
9		Benzaldehyde	0.894	0.788	11.9	149	0.00
10	C	Phenol	1.629	1.567	3.8	130	0.00
11		bis(2-Chloroethyl)ether	1.275	1.281	-0.5	144	0.00
12		1,3-Dichlorobenzene	1.595	1.579	1.0	142	0.00
13	C	1,4-Dichlorobenzene	1.692	1.656	2.1	141	0.00
14		1,2-Dichlorobenzene	1.559	1.541	1.2	139	0.00
15		Benzyl Alcohol	1.242	1.247	-0.4	138	0.00
16		2,2'-oxybis(1-Chloropropane	1.714	1.625	5.2	135	0.00
17		2-Methylphenol	1.146	1.131	1.3	136	0.00
18		Hexachloroethane	0.588	0.589	-0.2	140	0.00
19	P	n-Nitroso-di-n-propylamine	1.074	1.029	4.2	133	0.00
20		3+4-Methylphenols	1.517	1.413	6.9	131	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	139	0.00
22		Acetophenone	0.490	0.515	-5.1	137	0.00
23	S	Nitrobenzene-d5	0.361	0.373	-3.3	134	0.00
24		Nitrobenzene	0.368	0.379	-3.0	132	0.00
25		Isophorone	0.657	0.666	-1.4	131	0.00
26	C	2-Nitrophenol	0.172	0.182	-5.8	129	0.00
27		2,4-Dimethylphenol	0.284	0.288	-1.4	128	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.387	-3.5	131	0.00
29	C	2,4-Dichlorophenol	0.265	0.265	0.0	127	0.00
30		1,2,4-Trichlorobenzene	0.294	0.310	-5.4	133	0.00
31		Naphthalene	1.030	1.049	-1.8	132	0.00
32		Benzoic acid	0.158	0.132	16.5	107	0.03
33		4-Chloroaniline	0.416	0.421	-1.2	132	0.00
34	C	Hexachlorobutadiene	0.175	0.182	-4.0	134	0.00
35		Caprolactam	0.078	0.078	0.0	123	0.02
36	C	4-Chloro-3-methylphenol	0.303	0.309	-2.0	132	0.00
37		2-Methylnaphthalene	0.703	0.710	-1.0	133	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	134	0.00
39		1,2,4,5-Tetrachlorobenzene	0.494	0.495	-0.2	131	0.00
40	P	Hexachlorocyclopentadiene	0.202	0.241	-19.3	130	0.00
41	S	2,4,6-Tribromophenol	0.162	0.171	-5.6	130	0.00
42	C	2,4,6-Trichlorophenol	0.360	0.360	0.0	124	0.00
43		2,4,5-Trichlorophenol	0.367	0.351	4.4	119	0.00
44	S	2-Fluorobiphenyl	1.314	1.343	-2.2	131	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF020215\
Data File : BF077157.D
Acq On : 2 Feb 2015 14:09
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 03 09:57:22 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 03 00:09:26 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	1.483	1.524	-2.8	129	0.00
46		2-Chloronaphthalene	1.220	1.263	-3.5	134	0.00
47		2-Nitroaniline	0.370	0.380	-2.7	126	0.00
48		Acenaphthylene	2.058	2.060	-0.1	127	0.00
49		Dimethylphthalate	1.418	1.446	-2.0	132	0.00
50		2,6-Dinitrotoluene	0.283	0.308	-8.8	132	0.00
51	C	Acenaphthene	1.168	1.175	-0.6	129	0.00
52		3-Nitroaniline	0.339	0.348	-2.7	124	0.00
53	P	2,4-Dinitrophenol	0.095	0.110	-15.8	138	0.00
54		Dibenzofuran	1.701	1.717	-0.9	131	0.00
55	P	4-Nitrophenol	0.222	0.205	7.7	115	0.00
56		2,4-Dinitrotoluene	0.384	0.417	-8.6	131	0.00
57		Fluorene	1.441	1.449	-0.6	130	0.00
58		2,3,4,6-Tetrachlorophenol	0.268	0.267	0.4	123	0.00
59		Diethylphthalate	1.433	1.488	-3.8	133	0.00
60		4-Chlorophenyl-phenylether	0.590	0.607	-2.9	133	0.00
61		4-Nitroaniline	0.328	0.343	-4.6	127	0.00
62		Azobenzene	1.494	1.538	-2.9	132	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	140	0.00
64		4,6-Dinitro-2-methylphenol	0.104	0.126	-21.2	133	0.00
65	c	n-Nitrosodiphenylamine	0.713	0.705	1.1	127	0.00
66		4-Bromophenyl-phenylether	0.203	0.205	-1.0	129	0.00
67		Hexachlorobenzene	0.214	0.223	-4.2	139	0.00
68		Atrazine	0.216	0.217	-0.5	123	0.00
69	C	Pentachlorophenol	0.082	0.095	-15.9	146	0.00
70		Phenanthrene	1.247	1.225	1.8	131	0.00
71		Anthracene	1.216	1.246	-2.5	137	0.00
72		Carbazole	1.180	1.191	-0.9	132	0.00
73		Di-n-butylphthalate	1.365	1.467	-7.5	138	0.00
74	C	Fluoranthene	1.278	1.309	-2.4	132	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	147	0.00
76		Benzidine	0.640	0.689	-7.7	159#	0.00
77		Pyrene	1.371	1.359	0.9	133	0.00
78	S	Terphenyl-d14	0.869	0.859	1.2	135	0.00
79		Butylbenzylphthalate	0.621	0.659	-6.1	137	0.00
80		Benzo(a)anthracene	1.255	1.290	-2.8	140	0.00
81		3,3'-Dichlorobenzidine	0.399	0.426	-6.8	146	0.00
82		Chrysene	1.144	1.156	-1.0	137	0.00
83		Bis(2-ethylhexyl)phthalate	0.859	0.941	-9.5	149	0.00
84	c	Di-n-octyl phthalate	1.491	1.647	-10.5	148	0.00
85		Indeno(1,2,3-cd)pyrene	1.213	1.322	-9.0	146	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	144	-0.02
87		Benzo(b)fluoranthene	1.347	1.366	-1.4	153#	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF020215\
Data File : BF077157.D
Acq On : 2 Feb 2015 14:09
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 03 09:57:22 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 03 00:09:26 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	1.177	1.274	-8.2	137	0.00
89 C	Benzo(a)pyrene	1.188	1.236	-4.0	142	0.00
90	Dibenzo(a,h)anthracene	1.090	1.194	-9.5	149	-0.04
91	Benzo(g,h,i)perylene	1.084	1.147	-5.8	144	-0.03

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