Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF050815\
Data File : BF079108.D
Acq On : 10 May 2015 4:31
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 11 04:18:57 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 11 01:50:36 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	98	-0.05
2		1,4-Dioxane	0.505	0.503	0.4	98	-0.07
3		Pyridine	1.478	1.423	3.7	100	-0.08
4		n-Nitrosodimethylamine	0.633	0.637	-0.6	101	-0.08
5	S	2-Fluorophenol	1.162	1.192	-2.6	103	-0.05
6		Aniline	2.168	2.230	-2.9	102	-0.05
7	S	Phenol-d6	1.536	1.530	0.4	104	-0.05
8		2-Chlorophenol	1.318	1.345	-2.0	108	-0.06
9		Benzaldehyde	1.035	0.973	6.0	96	-0.05
10	C	Phenol	1.782	1.874	-5.2	105	-0.03
11		bis(2-Chloroethyl)ether	1.306	1.252	4.1	102	-0.06
12		1,3-Dichlorobenzene	1.481	1.516	-2.4	108	-0.06
13	C	1,4-Dichlorobenzene	1.524	1.522	0.1	104	-0.05
14		1,2-Dichlorobenzene	1.419	1.435	-1.1	104	-0.05
15		Benzyl Alcohol	1.140	1.112	2.5	101	-0.05
16		2,2'-oxybis(1-Chloropropane	1.998	1.999	-0.1	105	-0.05
17		2-Methylphenol	1.060	1.030	2.8	101	-0.05
18		Hexachloroethane	0.525	0.496	5.5	95	-0.06
19	P	n-Nitroso-di-n-propylamine	1.037	1.054	-1.6	100	-0.05
20		3+4-Methylphenols	1.413	1.424	-0.8	102	-0.05
21	I	Naphthalene-d8	1.000	1.000	0.0	96	-0.05
22		Acetophenone	0.452	0.449	0.7	102	-0.06
23	S	Nitrobenzene-d5	0.348	0.349	-0.3	101	-0.06
24		Nitrobenzene	0.345	0.334	3.2	101	-0.06
25		Isophorone	0.645	0.631	2.2	98	-0.06
26	C	2-Nitrophenol	0.143	0.149	-4.2	100	-0.05
27		2,4-Dimethylphenol	0.286	0.301	-5.2	103	-0.05
28		bis(2-Chloroethoxy)methane	0.398	0.402	-1.0	98	-0.05
29	C	2,4-Dichlorophenol	0.254	0.272	-7.1	107	-0.05
30		1,2,4-Trichlorobenzene	0.279	0.279	0.0	101	-0.05
31		Naphthalene	0.953	0.965	-1.3	104	-0.06
32		Benzoic acid	0.164	0.151	7.9	85	-0.06
33		4-Chloroaniline	0.403	0.396	1.7	101	-0.06
34	C	Hexachlorobutadiene	0.161	0.156	3.1	99	-0.06
35		Caprolactam	0.082	0.078	4.9	93	-0.07
36	C	4-Chloro-3-methylphenol	0.267	0.269	-0.7	94	-0.05
37		2-Methylnaphthalene	0.660	0.642	2.7	97	-0.06
38	I	Acenaphthene-d10	1.000	1.000	0.0	90	-0.06
39		1,2,4,5-Tetrachlorobenzene	0.563	0.577	-2.5	99	-0.06
40	P	Hexachlorocyclopentadiene	0.283	0.091	67.8#	30#	-0.06
41	S	2,4,6-Tribromophenol	0.216	0.225	-4.2	94	-0.06
42	C	2,4,6-Trichlorophenol	0.387	0.421	-8.8	100	-0.05
43		2,4,5-Trichlorophenol	0.392	0.423	-7.9	100	-0.05
44	S	2-Fluorobiphenyl	1.436	1.451	-1.0	95	-0.06

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF050815\
Data File : BF079108.D
Acq On : 10 May 2015 4:31
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 11 04:18:57 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 11 01:50:36 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	1.590	1.619	-1.8	99	-0.06
46		2-Chloronaphthalene	1.240	1.285	-3.6	102	-0.06
47		2-Nitroaniline	0.359	0.377	-5.0	96	-0.06
48		Acenaphthylene	2.036	2.152	-5.7	101	-0.06
49		Dimethylphthalate	1.468	1.519	-3.5	102	-0.06
50		2,6-Dinitrotoluene	0.290	0.283	2.4	91	-0.06
51	C	Acenaphthene	1.214	1.195	1.6	93	-0.06
52		3-Nitroaniline	0.362	0.390	-7.7	101	-0.06
53	P	2,4-Dinitrophenol	0.085	0.012#	85.9#	14#	-0.04
54		Dibenzofuran	1.754	1.736	1.0	94	-0.05
55	P	4-Nitrophenol	0.286	0.268	6.3	88	-0.03
56		2,4-Dinitrotoluene	0.383	0.381	0.5	89	-0.05
57		Fluorene	1.440	1.440	0.0	97	-0.06
58		2,3,4,6-Tetrachlorophenol	0.320	0.304	5.0	87	-0.05
59		Diethylphthalate	1.454	1.492	-2.6	98	-0.06
60		4-Chlorophenyl-phenylether	0.639	0.620	3.0	92	-0.06
61		4-Nitroaniline	0.362	0.385	-6.4	97	-0.06
62		Azobenzene	1.433	1.337	6.7	87	-0.06
63	I	Phenanthrene-d10	1.000	1.000	0.0	88	-0.05
64		4,6-Dinitro-2-methylphenol	0.080	0.018	77.5#	20#	-0.05
65	c	n-Nitrosodiphenylamine	0.666	0.662	0.6	89	-0.06
66		4-Bromophenyl-phenylether	0.208	0.207	0.5	94	-0.05
67		Hexachlorobenzene	0.238	0.236	0.8	94	-0.06
68		Atrazine	0.194	0.190	2.1	92	-0.06
69	C	Pentachlorophenol	0.120	0.108	10.0	80	-0.04
70		Phenanthrene	1.119	1.062	5.1	87	-0.06
71		Anthracene	1.098	1.125	-2.5	95	-0.06
72		Carbazole	1.046	1.083	-3.5	95	-0.05
73		Di-n-butylphthalate	1.255	1.330	-6.0	93	-0.06
74	C	Fluoranthene	1.166	1.178	-1.0	92	-0.06
75	I	Chrysene-d12	1.000	1.000	0.0	80	-0.08
76		Benzidine	0.539	0.681	-26.3#	98	-0.06
77		Pyrene	1.152	1.200	-4.2	89	-0.06
78	S	Terphenyl-d14	0.835	0.834	0.1	85	-0.06
79		Butylbenzylphthalate	0.504	0.580	-15.1	90	-0.06
80		Benzo(a)anthracene	1.095	1.127	-2.9	87	-0.08
81		3,3'-Dichlorobenzidine	0.390	0.436	-11.8	90	-0.07
82		Chrysene	1.053	1.038	1.4	86	-0.08
83		Bis(2-ethylhexyl)phthalate	0.763	0.841	-10.2	86	-0.08
84	c	Di-n-octyl phthalate	1.344	1.521	-13.2	95	-0.13
85		Indeno(1,2,3-cd)pyrene	1.342	1.336	0.4	84	-0.22
86	I	Perylene-d12	1.000	1.000	0.0	85	-0.17
87		Benzo(b)fluoranthene	1.095	1.128	-3.0	90	-0.15

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF050815\
Data File : BF079108.D
Acq On : 10 May 2015 4:31
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 11 04:18:57 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon May 11 01:50:36 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	1.060	1.039	2.0	85	-0.16
89 C	Benzo(a)pyrene	1.008	1.022	-1.4	89	-0.17
90	Dibenzo(a,h)anthracene	1.071	1.020	4.8	83	-0.23
91	Benzo(g,h,i)perylene	1.074	1.014	5.6	84	-0.24

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