Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079422.D
Acq On : 22 May 2015 2:23
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 22 05:02:09 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 22 05:01:56 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	58	-0.06
2		1,4-Dioxane	0.505	0.625	-23.8	73	-0.08
3		Pyridine	1.478	1.526	-3.2	64	-0.09
4		n-Nitrosodimethylamine	0.633	0.591	6.6	56	-0.08
5	S	2-Fluorophenol	1.162	1.237	-6.5	64	-0.07
6		Aniline	2.168	2.205	-1.7	61	-0.06
7	S	Phenol-d6	1.536	1.603	-4.4	65	-0.06
8		2-Chlorophenol	1.318	1.436	-9.0	69	-0.07
9		Benzaldehyde	1.035	0.972	6.1	57	-0.06
10	C	Phenol	1.782	1.900	-6.6	64	-0.06
11		bis(2-Chloroethyl)ether	1.306	1.381	-5.7	67	-0.06
12		1,3-Dichlorobenzene	1.481	1.578	-6.5	67	-0.06
13	C	1,4-Dichlorobenzene	1.524	1.615	-6.0	66	-0.06
14		1,2-Dichlorobenzene	1.419	1.522	-7.3	66	-0.06
15		Benzyl Alcohol	1.140	1.085	4.8	59	-0.06
16		2,2'-oxybis(1-Chloropropane	1.998	2.001	-0.2	63	-0.06
17		2-Methylphenol	1.060	1.147	-8.2	67	-0.06
18		Hexachloroethane	0.525	0.563	-7.2	64	-0.06
19	P	n-Nitroso-di-n-propylamine	1.037	1.071	-3.3	61	-0.06
20		3+4-Methylphenols	1.413	1.507	-6.7	64	-0.06
21	I	Naphthalene-d8	1.000	1.000	0.0	59	-0.06
22		Acetophenone	0.452	0.475	-5.1	67	-0.06
23	S	Nitrobenzene-d5	0.348	0.367	-5.5	65	-0.06
24		Nitrobenzene	0.345	0.364	-5.5	68	-0.06
25		Isophorone	0.645	0.676	-4.8	64	-0.06
26	C	2-Nitrophenol	0.143	0.146	-2.1	60	-0.06
27		2,4-Dimethylphenol	0.286	0.288	-0.7	61	-0.05
28		bis(2-Chloroethoxy)methane	0.398	0.425	-6.8	64	-0.06
29	C	2,4-Dichlorophenol	0.254	0.276	-8.7	67	-0.06
30		1,2,4-Trichlorobenzene	0.279	0.301	-7.9	67	-0.06
31		Naphthalene	0.953	1.045	-9.7	69	-0.06
32		Benzoic acid	0.164	0.180	-9.8	62	-0.07
33		4-Chloroaniline	0.403	0.435	-7.9	68	-0.06
34	C	Hexachlorobutadiene	0.161	0.169	-5.0	67	-0.06
35		Caprolactam	0.082	0.087	-6.1	64	-0.07
36	C	4-Chloro-3-methylphenol	0.267	0.300	-12.4	65	-0.06
37		2-Methylnaphthalene	0.660	0.717	-8.6	67	-0.06
38	I	Acenaphthene-d10	1.000	1.000	0.0	58	-0.07
39		1,2,4,5-Tetrachlorobenzene	0.563	0.598	-6.2	67	-0.07
40	P	Hexachlorocyclopentadiene	0.283	0.319	-12.7	68	-0.06
41	S	2,4,6-Tribromophenol	0.216	0.233	-7.9	63	-0.07
42	C	2,4,6-Trichlorophenol	0.387	0.404	-4.4	62	-0.07
43		2,4,5-Trichlorophenol	0.392	0.401	-2.3	62	-0.06
44	S	2-Fluorobiphenyl	1.436	1.582	-10.2	67	-0.06
45		1,1'-Biphenyl	1.590	1.705	-7.2	67	-0.06
46		2-Chloronaphthalene	1.240	1.347	-8.6	69	-0.06
47		2-Nitroaniline	0.359	0.375	-4.5	62	-0.06
48		Acenaphthylene	2.036	2.263	-11.1	69	-0.07

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF052115\
Data File : BF079422.D
Acq On : 22 May 2015 2:23
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 22 05:02:09 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF043015.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 22 05:01:56 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)
49		Dimethylphthalate	1.468	1.522	-3.7	66	-0.07
50		2,6-Dinitrotoluene	0.290	0.321	-10.7	67	-0.06
51	C	Acenaphthene	1.214	1.277	-5.2	64	-0.06
52		3-Nitroaniline	0.362	0.396	-9.4	67	-0.07
53	P	2,4-Dinitrophenol	0.085	0.045#	47.1#	32#	-0.06
54		Dibenzofuran	1.754	1.887	-7.6	66	-0.06
55	P	4-Nitrophenol	0.286	0.208	27.3#	45#	-0.05
56		2,4-Dinitrotoluene	0.383	0.413	-7.8	62	-0.07
57		Fluorene	1.440	1.568	-8.9	69	-0.06
58		2,3,4,6-Tetrachlorophenol	0.320	0.308	3.8	57	-0.06
59		Diethylphthalate	1.454	1.564	-7.6	67	-0.06
60		4-Chlorophenyl-phenylether	0.639	0.681	-6.6	66	-0.06
61		4-Nitroaniline	0.362	0.405	-11.9	66	-0.07
62		Azobenzene	1.433	1.508	-5.2	63	-0.06
63	I	Phenanthrene-d10	1.000	1.000	0.0	59	-0.07
64		4,6-Dinitro-2-methylphenol	0.080	0.061	23.8	47#	-0.07
65	c	n-Nitrosodiphenylamine	0.666	0.709	-6.5	65	-0.06
66		4-Bromophenyl-phenylether	0.208	0.219	-5.3	67	-0.06
67		Hexachlorobenzene	0.238	0.248	-4.2	67	-0.07
68		Atrazine	0.194	0.193	0.5	63	-0.07
69	C	Pentachlorophenol	0.120	0.090	25.0#	45#	-0.07
70		Phenanthrene	1.119	1.174	-4.9	65	-0.06
71		Anthracene	1.098	1.151	-4.8	66	-0.07
72		Carbazole	1.046	1.123	-7.4	67	-0.06
73		Di-n-butylphthalate	1.255	1.309	-4.3	62	-0.07
74	C	Fluoranthene	1.166	1.248	-7.0	66	-0.07
75	I	Chrysene-d12	1.000	1.000	0.0	55	-0.09
76		Benzidine	0.539	0.582	-8.0	58	-0.07
77		Pyrene	1.152	1.296	-12.5	66	-0.07
78	S	Terphenyl-d14	0.835	0.933	-11.7	65	-0.07
79		Butylbenzylphthalate	0.504	0.569	-12.9	61	-0.07
80		Benzo(a)anthracene	1.095	1.178	-7.6	62	-0.09
81		3,3'-Dichlorobenzidine	0.390	0.422	-8.2	60	-0.08
82		Chrysene	1.053	1.143	-8.5	65	-0.09
83		Bis(2-ethylhexyl)phthalate	0.763	0.870	-14.0	61	-0.08
84	c	Di-n-octyl phthalate	1.344	1.513	-12.6	65	-0.13
85		Indeno(1,2,3-cd)pyrene	1.342	1.548	-15.4	67	-0.26
86	I	Perylene-d12	1.000	1.000	0.0	59	-0.19
87		Benzo(b)fluoranthene	1.095	1.433	-30.9#	79	-0.16
88		Benzo(k)fluoranthene	1.060	0.910	14.2	52	-0.16
89	C	Benzo(a)pyrene	1.008	1.119	-11.0	67	-0.19
90		Dibenzo(a,h)anthracene	1.071	1.218	-13.7	68	-0.27
91		Benzo(g,h,i)perylene	1.074	1.209	-12.6	69	-0.27

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