Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF081117\
Data File : BF097620.D
Acq On : 12 Aug 2017 4:58
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 14 11:07:17 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 11 15:55:57 2017
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	69	0.00
2		1,4-Dioxane	0.546	0.559	-2.4	76	-0.05
3		Pyridine	1.593	1.599	-0.4	64	-0.04
4		n-Nitrosodimethylamine	0.924	0.911	1.4	66	-0.05
5	S	2-Fluorophenol	1.258	1.316	-4.6	67	0.02
6		Aniline	2.243	2.182	2.7	70	0.00
7	S	Phenol-d6	1.561	1.523	2.4	69	0.02
8		2-Chlorophenol	1.442	1.401	2.8	68	0.02
9		Benzaldehyde	1.029	1.077	-4.7	73	0.00
10	C	Phenol	1.709	1.775	-3.9	67	0.03
11		bis(2-Chloroethyl)ether	1.371	1.409	-2.8	72	0.00
12		1,3-Dichlorobenzene	1.615	1.544	4.4	67	0.00
13	C	1,4-Dichlorobenzene	1.644	1.723	-4.8	74	0.00
14		1,2-Dichlorobenzene	1.500	1.625	-8.3	77	0.00
15		Benzyl Alcohol	1.079	1.207	-11.9	75	0.01
16		2,2'-oxybis(1-Chloropropane	3.078	3.314	-7.7	77	0.02
17		2-Methylphenol	1.118	1.187	-6.2	76	0.03
18		Hexachloroethane	0.563	0.590	-4.8	75	0.00
19	P	n-Nitroso-di-n-propylamine	1.086	1.156	-6.4	73	0.00
20		3+4-Methylphenols	1.369	1.509	-10.2	75	0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	62	0.00
22		Acetophenone	0.472	0.564	-19.5	73	0.00
23	S	Nitrobenzene-d5	0.319	0.387	-21.3	73	0.00
24		Nitrobenzene	0.348	0.433	-24.4	74	0.00
25		Isophorone	0.588	0.673	-14.5	68	0.00
26	C	2-Nitrophenol	0.176	0.214	-21.6#	71	0.00
27		2,4-Dimethylphenol	0.293	0.331	-13.0	68	0.01
28		bis(2-Chloroethoxy)methane	0.425	0.496	-16.7	70	0.00
29	C	2,4-Dichlorophenol	0.259	0.277	-6.9	65	0.01
30		1,2,4-Trichlorobenzene	0.256	0.297	-16.0	71	0.00
31		Naphthalene	1.002	1.094	-9.2	68	0.00
32		Benzoic acid	0.158	0.128	19.0	44#	-0.01
33		4-Chloroaniline	0.415	0.511	-23.1	74	0.00
34	C	Hexachlorobutadiene	0.140	0.171	-22.1#	76	0.00
35		Caprolactam	0.083	0.103	-24.1	73	0.00
36	C	4-Chloro-3-methylphenol	0.290	0.346	-19.3	69	0.02
37		2-Methylnaphthalene	0.630	0.751	-19.2	73	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	66	0.00
39		1,2,4,5-Tetrachlorobenzene	0.566	0.643	-13.6	75	0.00
40	P	Hexachlorocyclopentadiene	0.072	0.008#	88.9#	7#	0.00
41	S	2,4,6-Tribromophenol	0.152	0.149	2.0	65	0.00
42	C	2,4,6-Trichlorophenol	0.373	0.430	-15.3	73	0.00
43		2,4,5-Trichlorophenol	0.352	0.389	-10.5	67	0.02
44	S	2-Fluorobiphenyl	1.325	1.414	-6.7	72	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF081117\
Data File : BF097620.D
Acq On : 12 Aug 2017 4:58
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 14 11:07:17 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 11 15:55:57 2017
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	1.716	1.889	-10.1	74	0.00
46		2-Chloronaphthalene	1.317	1.422	-8.0	72	0.00
47		2-Nitroaniline	0.516	0.536	-3.9	62	0.01
48		Acenaphthylene	2.094	2.151	-2.7	68	0.00
49		Dimethylphthalate	1.541	1.606	-4.2	69	0.00
50		2,6-Dinitrotoluene	0.318	0.338	-6.3	68	0.00
51	C	Acenaphthene	1.255	1.259	-0.3	68	0.00
52		3-Nitroaniline	0.417	0.431	-3.4	67	0.00
53	P	2,4-Dinitrophenol	0.131	0.024#	81.7#	13#	0.04
54		Dibenzofuran	1.757	1.790	-1.9	68	0.00
55	P	4-Nitrophenol	0.159	0.054	66.0#	23#	0.06
56		2,4-Dinitrotoluene	0.425	0.448	-5.4	67	0.00
57		Fluorene	1.445	1.500	-3.8	69	0.00
58		2,3,4,6-Tetrachlorophenol	0.242	0.207	14.5	53	0.02
59		Diethylphthalate	1.456	1.567	-7.6	72	0.00
60		4-Chlorophenyl-phenylether	0.617	0.652	-5.7	71	0.00
61		4-Nitroaniline	0.398	0.383	3.8	60	0.00
62		Azobenzene	1.448	1.434	1.0	65	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	63	0.00
64		4,6-Dinitro-2-methylphenol	0.122	0.084	31.1#	41#	0.00
65	c	n-Nitrosodiphenylamine	0.739	0.783	-6.0	69	0.00
66		4-Bromophenyl-phenylether	0.207	0.228	-10.1	71	0.00
67		Hexachlorobenzene	0.226	0.239	-5.8	68	0.00
68		Atrazine	0.185	0.203	-9.7	69	0.01
69	C	Pentachlorophenol	0.037	0.002	94.6#	3#	0.04
70		Phenanthrene	1.152	1.162	-0.9	66	0.00
71		Anthracene	1.179	1.192	-1.1	66	0.00
72		Carbazole	1.122	1.113	0.8	65	0.01
73		Di-n-butylphthalate	1.313	1.460	-11.2	71	0.00
74	C	Fluoranthene	1.041	1.103	-6.0	64	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	64	0.00
76		Benzidine	0.659	0.567	14.0	55	0.00
77		Pyrene	1.654	1.515	8.4	64	0.00
78	S	Terphenyl-d14	0.966	0.904	6.4	69	0.00
79		Butylbenzylphthalate	0.691	0.746	-8.0	66	0.00
80		Benzo(a)anthracene	1.168	1.179	-0.9	66	0.00
81		3,3'-Dichlorobenzidine	0.403	0.436	-8.2	69	0.00
82		Chrysene	1.162	1.172	-0.9	65	0.00
83		Bis(2-ethylhexyl)phthalate	0.983	1.074	-9.3	67	0.00
84	c	Di-n-octyl phthalate	1.717	1.764	-2.7	68	0.00
85		Indeno(1,2,3-cd)pyrene	0.904	1.039	-14.9	94	0.01
86	I	Perylene-d12	1.000	1.000	0.0	76	0.00
87		Benzo(b)fluoranthene	1.201	1.140	5.1	75	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF081117\
Data File : BF097620.D
Acq On : 12 Aug 2017 4:58
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 14 11:07:17 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 11 15:55:57 2017
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	1.210	1.204	0.5	75	0.00
89 C	Benzo(a)pyrene	1.102	1.128	-2.4	80	0.01
90	Dibenzo(a,h)anthracene	0.789	0.939	-19.0	94	0.02
91	Benzo(g,h,i)perylene	0.866	0.932	-7.6	89	0.02

(#) = Out of Range SPCC's out = 2 CCC's out = 3