Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051617\
Data File : BF095158.D
Acq On : 16 May 2017 15:56
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 17 13:06:08 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 03 17:24:51 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	96	0.05
2		1,4-Dioxane	0.588	0.519	11.7	90	0.15
3		Pyridine	1.364	1.300	4.7	94	0.19
4		n-Nitrosodimethylamine	0.568	0.478	15.8	84	0.18
5	S	2-Fluorophenol	1.200	1.178	1.8	95	0.06
6		Aniline	1.817	1.882	-3.6	96	0.05
7	S	Phenol-d6	1.439	1.463	-1.7	98	0.04
8		2-Chlorophenol	1.365	1.405	-2.9	101	0.05
9		Benzaldehyde	0.879	0.779	11.4	84	0.06
10	C	Phenol	1.517	1.434	5.5	92	0.04
11		bis(2-Chloroethyl)ether	1.252	1.284	-2.6	99	0.05
12		1,3-Dichlorobenzene	1.549	1.600	-3.3	107	0.05
13	C	1,4-Dichlorobenzene	1.571	1.624	-3.4	99	0.04
14		1,2-Dichlorobenzene	1.466	1.517	-3.5	101	0.04
15		Benzyl Alcohol	0.926	1.016	-9.7	102	0.05
16		2,2'-oxybis(1-Chloropropane	1.772	1.815	-2.4	102	0.04
17		2-Methylphenol	1.117	1.149	-2.9	103	0.04
18		Hexachloroethane	0.532	0.549	-3.2	99	0.04
19	P	n-Nitroso-di-n-propylamine	0.973	1.004	-3.2	102	0.04
20		3+4-Methylphenols	1.518	1.563	-3.0	100	0.04
21	I	Naphthalene-d8	1.000	1.000	0.0	98	0.04
22		Acetophenone	0.480	0.478	0.4	100	0.05
23	S	Nitrobenzene-d5	0.317	0.317	0.0	98	0.05
24		Nitrobenzene	0.332	0.330	0.6	101	0.05
25		Isophorone	0.566	0.587	-3.7	103	0.05
26	C	2-Nitrophenol	0.179	0.189	-5.6	102	0.05
27		2,4-Dimethylphenol	0.290	0.289	0.3	102	0.04
28		bis(2-Chloroethoxy)methane	0.392	0.359	8.4	91	0.04
29	C	2,4-Dichlorophenol	0.274	0.260	5.1	98	0.05
30		1,2,4-Trichlorobenzene	0.293	0.296	-1.0	102	0.04
31		Naphthalene	1.005	0.989	1.6	100	0.04
32		Benzoic acid	0.177	0.131	26.0#	76	0.05
33		4-Chloroaniline	0.375	0.387	-3.2	103	0.04
34	C	Hexachlorobutadiene	0.155	0.146	5.8	95	0.03
35		Caprolactam	0.082	0.076	7.3	88	0.06
36	C	4-Chloro-3-methylphenol	0.293	0.293	0.0	100	0.04
37		2-Methylnaphthalene	0.660	0.664	-0.6	102	0.04
38	I	Acenaphthene-d10	1.000	1.000	0.0	98	0.04
39		1,2,4,5-Tetrachlorobenzene	0.615	0.647	-5.2	98	0.04
40	P	Hexachlorocyclopentadiene	0.221	0.262	-18.6	106	0.03
41	S	2,4,6-Tribromophenol	0.164	0.171	-4.3	95	0.04
42	C	2,4,6-Trichlorophenol	0.411	0.394	4.1	89	0.04
43		2,4,5-Trichlorophenol	0.441	0.416	5.7	87	0.05
44	S	2-Fluorobiphenyl	1.390	1.324	4.7	91	0.04

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051617\
Data File : BF095158.D
Acq On : 16 May 2017 15:56
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 17 13:06:08 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 03 17:24:51 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	1.684	1.741	-3.4	96	0.04
46		2-Chloronaphthalene	1.360	1.390	-2.2	97	0.04
47		2-Nitroaniline	0.372	0.387	-4.0	99	0.05
48		Acenaphthylene	2.130	2.252	-5.7	101	0.04
49		Dimethylphthalate	1.642	1.748	-6.5	101	0.04
50		2,6-Dinitrotoluene	0.335	0.365	-9.0	102	0.05
51	C	Acenaphthene	1.272	1.267	0.4	95	0.04
52		3-Nitroaniline	0.360	0.359	0.3	93	0.06
53	P	2,4-Dinitrophenol	0.156	0.142	9.0	83	0.07
54		Dibenzofuran	1.760	1.888	-7.3	104	0.04
55	P	4-Nitrophenol	0.216	0.187	13.4	77	0.08
56		2,4-Dinitrotoluene	0.430	0.474	-10.2	102	0.06
57		Fluorene	1.531	1.428	6.7	91	0.04
58		2,3,4,6-Tetrachlorophenol	0.278	0.309	-11.2	106	0.05
59		Diethylphthalate	1.574	1.517	3.6	95	0.04
60		4-Chlorophenyl-phenylether	0.673	0.659	2.1	93	0.04
61		4-Nitroaniline	0.330	0.268	18.8	78	0.05
62		Azobenzene	1.265	1.257	0.6	92	0.04
63	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.04
64		4,6-Dinitro-2-methylphenol	0.134	0.128	4.5	90	0.06
65	c	n-Nitrosodiphenylamine	0.726	0.694	4.4	93	0.04
66		4-Bromophenyl-phenylether	0.211	0.226	-7.1	100	0.04
67		Hexachlorobenzene	0.217	0.231	-6.5	103	0.05
68		Atrazine	0.205	0.201	2.0	99	0.04
69	C	Pentachlorophenol	0.085	0.076	10.6	91	0.05
70		Phenanthrene	1.149	1.145	0.3	98	0.04
71		Anthracene	1.174	1.220	-3.9	101	0.04
72		Carbazole	1.068	1.082	-1.3	100	0.05
73		Di-n-butylphthalate	1.332	1.329	0.2	95	0.03
74	C	Fluoranthene	1.179	1.094	7.2	89	0.04
75	I	Chrysene-d12	1.000	1.000	0.0	95	0.04
76		Benzidine	0.465	0.398	14.4	71	0.05
77		Pyrene	1.496	1.456	2.7	91	0.04
78	S	Terphenyl-d14	0.908	0.891	1.9	94	0.04
79		Butylbenzylphthalate	0.696	0.719	-3.3	96	0.04
80		Benzo(a)anthracene	1.164	1.152	1.0	94	0.05
81		3,3'-Dichlorobenzidine	0.402	0.381	5.2	86	0.04
82		Chrysene	1.125	1.145	-1.8	96	0.04
83		Bis(2-ethylhexyl)phthalate	0.991	1.108	-11.8	103	0.03
84	c	Di-n-octyl phthalate	1.625	1.884	-15.9	107	0.03
85		Indeno(1,2,3-cd)pyrene	0.914	0.769	15.9	81	0.09
86	I	Perylene-d12	1.000	1.000	0.0	92	0.05
87		Benzo(b)fluoranthene	1.236	1.188	3.9	81	0.05

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF051617\
Data File : BF095158.D
Acq On : 16 May 2017 15:56
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 17 13:06:08 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF050217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 03 17:24:51 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	1.192	1.310	-9.9	98	0.05
89 C	Benzo(a)pyrene	1.140	1.110	2.6	85	0.05
90	Dibenzo(a,h)anthracene	0.942	0.829	12.0	79	0.08
91	Benzo(g,h,i)perylene	0.924	0.801	13.3	80	0.10

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