Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF060717\
Data File : BF095744.D
Acq On : 7 Jun 2017 14:05
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 08 04:04:31 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 05 16:15:31 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	71	0.02
2		1,4-Dioxane	0.597	0.563	5.7	65	0.00
3		Pyridine	1.403	1.347	4.0	66	0.01
4		n-Nitrosodimethylamine	0.534	0.520	2.6	67	0.00
5	S	2-Fluorophenol	1.255	1.304	-3.9	68	0.01
6		Aniline	1.966	1.998	-1.6	68	0.02
7	S	Phenol-d6	1.503	1.558	-3.7	68	0.01
8		2-Chlorophenol	1.335	1.398	-4.7	69	0.01
9		Benzaldehyde	1.014	1.010	0.4	67	0.01
10	C	Phenol	1.559	1.639	-5.1	67	0.01
11		bis(2-Chloroethyl)ether	1.235	1.315	-6.5	70	0.01
12		1,3-Dichlorobenzene	1.511	1.548	-2.4	72	0.02
13	C	1,4-Dichlorobenzene	1.532	1.597	-4.2	72	0.01
14		1,2-Dichlorobenzene	1.412	1.504	-6.5	73	0.01
15		Benzyl Alcohol	1.031	1.089	-5.6	71	0.01
16		2,2'-oxybis(1-Chloropropane	1.735	1.767	-1.8	70	0.02
17		2-Methylphenol	1.120	1.149	-2.6	70	0.01
18		Hexachloroethane	0.506	0.531	-4.9	72	0.01
19	P	n-Nitroso-di-n-propylamine	0.952	1.006	-5.7	72	0.02
20		3+4-Methylphenols	1.422	1.529	-7.5	73	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	71	0.02
22		Acetophenone	0.468	0.479	-2.4	72	0.02
23	S	Nitrobenzene-d5	0.322	0.327	-1.6	71	0.02
24		Nitrobenzene	0.344	0.347	-0.9	71	0.01
25		Isophorone	0.601	0.604	-0.5	71	0.02
26	C	2-Nitrophenol	0.180	0.192	-6.7	73	0.02
27		2,4-Dimethylphenol	0.280	0.288	-2.9	73	0.01
28		bis(2-Chloroethoxy)methane	0.396	0.405	-2.3	71	0.02
29	C	2,4-Dichlorophenol	0.269	0.284	-5.6	74	0.01
30		1,2,4-Trichlorobenzene	0.280	0.293	-4.6	74	0.01
31		Naphthalene	1.004	1.021	-1.7	73	0.02
32		Benzoic acid	0.161	0.166	-3.1	69	0.01
33		4-Chloroaniline	0.392	0.409	-4.3	74	0.02
34	C	Hexachlorobutadiene	0.145	0.153	-5.5	75	0.01
35		Caprolactam	0.080	0.085	-6.3	71	0.02
36	C	4-Chloro-3-methylphenol	0.282	0.290	-2.8	72	0.00
37		2-Methylnaphthalene	0.678	0.690	-1.8	73	0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	78	0.02
39		1,2,4,5-Tetrachlorobenzene	0.599	0.608	-1.5	75	0.02
40	P	Hexachlorocyclopentadiene	0.138	0.130	5.8	67	0.01
41	S	2,4,6-Tribromophenol	0.177	0.182	-2.8	78	0.01
42	C	2,4,6-Trichlorophenol	0.385	0.396	-2.9	74	0.01
43		2,4,5-Trichlorophenol	0.409	0.410	-0.2	70	0.01
44	S	2-Fluorobiphenyl	1.437	1.440	-0.2	76	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF060717\
Data File : BF095744.D
Acq On : 7 Jun 2017 14:05
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 08 04:04:31 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 05 16:15:31 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	1.648	1.646	0.1	74	0.02
46		2-Chloronaphthalene	1.288	1.283	0.4	74	0.01
47		2-Nitroaniline	0.377	0.372	1.3	68	0.02
48		Acenaphthylene	2.202	2.180	1.0	72	0.02
49		Dimethylphthalate	1.519	1.504	1.0	73	0.01
50		2,6-Dinitrotoluene	0.331	0.332	-0.3	70	0.01
51	C	Acenaphthene	1.272	1.268	0.3	77	0.01
52		3-Nitroaniline	0.386	0.382	1.0	76	0.02
53	P	2,4-Dinitrophenol	0.160	0.165	-3.1	77	0.01
54		Dibenzofuran	1.790	1.792	-0.1	78	0.02
55	P	4-Nitrophenol	0.201	0.189	6.0	68	0.01
56		2,4-Dinitrotoluene	0.447	0.455	-1.8	76	0.01
57		Fluorene	1.558	1.561	-0.2	78	0.02
58		2,3,4,6-Tetrachlorophenol	0.280	0.290	-3.6	77	0.01
59		Diethylphthalate	1.461	1.460	0.1	77	0.02
60		4-Chlorophenyl-phenylether	0.677	0.680	-0.4	77	0.02
61		4-Nitroaniline	0.360	0.360	0.0	75	0.02
62		Azobenzene	1.324	1.279	3.4	75	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	75	0.01
64		4,6-Dinitro-2-methylphenol	0.131	0.137	-4.6	75	0.01
65	c	n-Nitrosodiphenylamine	0.719	0.718	0.1	77	0.02
66		4-Bromophenyl-phenylether	0.208	0.208	0.0	75	0.01
67		Hexachlorobenzene	0.205	0.213	-3.9	78	0.01
68		Atrazine	0.200	0.206	-3.0	79	0.01
69	C	Pentachlorophenol	0.067	0.070	-4.5	77	0.01
70		Phenanthrene	1.154	1.164	-0.9	78	0.01
71		Anthracene	1.205	1.221	-1.3	79	0.02
72		Carbazole	1.174	1.201	-2.3	77	0.02
73		Di-n-butylphthalate	1.249	1.322	-5.8	79	0.02
74	C	Fluoranthene	1.142	1.176	-3.0	78	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	78	0.01
76		Benzidine	0.768	0.728	5.2	72	0.01
77		Pyrene	1.492	1.521	-1.9	79	0.01
78	S	Terphenyl-d14	0.806	0.825	-2.4	80	0.01
79		Butylbenzylphthalate	0.652	0.670	-2.8	77	0.01
80		Benzo(a)anthracene	1.163	1.194	-2.7	78	0.01
81		3,3'-Dichlorobenzidine	0.413	0.433	-4.8	79	0.01
82		Chrysene	1.162	1.156	0.5	76	0.01
83		Bis(2-ethylhexyl)phthalate	0.919	0.968	-5.3	79	0.01
84	c	Di-n-octyl phthalate	1.515	1.570	-3.6	78	0.01
85		Indeno(1,2,3-cd)pyrene	0.994	0.904	9.1	75	0.01
86	I	Perylene-d12	1.000	1.000	0.0	73	0.01
87		Benzo(b)fluoranthene	1.252	1.340	-7.0	74	0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF060717\
Data File : BF095744.D
Acq On : 7 Jun 2017 14:05
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 08 04:04:31 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 05 16:15:31 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	1.197	1.183	1.2	74	0.01
89 C	Benzo(a)pyrene	1.151	1.163	-1.0	74	0.01
90	Dibenzo(a,h)anthracene	0.976	0.933	4.4	74	0.02
91	Benzo(g,h,i)perylene	0.995	0.940	5.5	73	0.02

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