Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
Data File : BF092141.D
Acq On : 9 Jan 2017 17:00
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 09 22:40:52 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 06 15:43:51 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	76	-0.01
2		1,4-Dioxane	0.590	0.693	-17.5	88	0.02
3		Pyridine	2.058	1.772	13.9	63	0.01
4		n-Nitrosodimethylamine	0.776	0.725	6.6	67	0.01
5	S	2-Fluorophenol	1.198	1.270	-6.0	83	0.00
6		Aniline	1.899	1.720	9.4	68	-0.01
7	S	Phenol-d6	1.462	1.391	4.9	70	0.00
8		2-Chlorophenol	1.362	1.321	3.0	72	-0.01
9		Benzaldehyde	0.904	0.735	18.7	66	-0.01
10	C	Phenol	1.666	1.704	-2.3	71	0.00
11		bis(2-Chloroethyl)ether	1.326	1.289	2.8	68	-0.01
12		1,3-Dichlorobenzene	1.455	1.530	-5.2	75	-0.01
13	C	1,4-Dichlorobenzene	1.480	1.561	-5.5	78	-0.01
14		1,2-Dichlorobenzene	1.285	1.341	-4.4	80	-0.01
15		Benzyl Alcohol	1.136	1.148	-1.1	75	-0.01
16		2,2'-oxybis(1-Chloropropane	1.975	2.177	-10.2	83	-0.02
17		2-Methylphenol	1.096	1.193	-8.9	82	0.00
18		Hexachloroethane	0.549	0.490	10.7	64	-0.01
19	P	n-Nitroso-di-n-propylamine	0.973	0.925	4.9	68	-0.01
20		3+4-Methylphenols	1.307	1.409	-7.8	77	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	67	-0.01
22		Acetophenone	0.493	0.545	-10.5	73	-0.01
23	S	Nitrobenzene-d5	0.360	0.371	-3.1	72	0.00
24		Nitrobenzene	0.383	0.348	9.1	55	-0.01
25		Isophorone	0.664	0.644	3.0	65	-0.01
26	C	2-Nitrophenol	0.189	0.200	-5.8	72	-0.01
27		2,4-Dimethylphenol	0.313	0.284	9.3	56	-0.01
28		bis(2-Chloroethoxy)methane	0.402	0.358	10.9	58	-0.01
29	C	2,4-Dichlorophenol	0.265	0.267	-0.8	65	-0.01
30		1,2,4-Trichlorobenzene	0.291	0.278	4.5	61	-0.01
31		Naphthalene	0.915	0.971	-6.1	65	-0.01
32		Benzoic acid	0.232	0.240	-3.4	65	0.01
33		4-Chloroaniline	0.413	0.379	8.2	60	-0.01
34	C	Hexachlorobutadiene	0.153	0.162	-5.9	69	-0.01
35		Caprolactam	0.087	0.085	2.3	63	0.00
36	C	4-Chloro-3-methylphenol	0.305	0.289	5.2	59	0.00
37		2-Methylnaphthalene	0.628	0.596	5.1	69	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	66	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.505	0.508	-0.6	59	-0.01
40	P	Hexachlorocyclopentadiene	0.199	0.204	-2.5	58	-0.01
41	S	2,4,6-Tribromophenol	0.166	0.177	-6.6	72	-0.01
42	C	2,4,6-Trichlorophenol	0.353	0.375	-6.2	63	0.00
43		2,4,5-Trichlorophenol	0.364	0.375	-3.0	65	0.00
44	S	2-Fluorobiphenyl	1.042	1.226	-17.7	80	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
Data File : BF092141.D
Acq On : 9 Jan 2017 17:00
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 09 22:40:52 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 06 15:43:51 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	1.369	1.405	-2.6	59	-0.01
46		2-Chloronaphthalene	1.084	1.098	-1.3	63	-0.01
47		2-Nitroaniline	0.386	0.393	-1.8	59	-0.01
48		Acenaphthylene	1.668	1.677	-0.5	65	-0.01
49		Dimethylphthalate	1.279	1.407	-10.0	68	0.00
50		2,6-Dinitrotoluene	0.280	0.332	-18.6	77	0.00
51	C	Acenaphthene	1.131	1.057	6.5	61	-0.01
52		3-Nitroaniline	0.346	0.417	-20.5	69	0.00
53	P	2,4-Dinitrophenol	0.156	0.230	-47.4#	84	-0.01
54		Dibenzofuran	1.527	1.382	9.5	64	-0.01
55	P	4-Nitrophenol	0.235	0.250	-6.4	63	0.01
56		2,4-Dinitrotoluene	0.351	0.358	-2.0	68	-0.01
57		Fluorene	1.111	1.161	-4.5	65	-0.01
58		2,3,4,6-Tetrachlorophenol	0.288	0.291	-1.0	61	0.00
59		Diethylphthalate	1.242	1.287	-3.6	62	-0.01
60		4-Chlorophenyl-phenylether	0.486	0.551	-13.4	74	0.00
61		4-Nitroaniline	0.359	0.364	-1.4	59	-0.01
62		Azobenzene	1.368	1.485	-8.6	72	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	67	-0.01
64		4,6-Dinitro-2-methylphenol	0.121	0.146	-20.7	75	0.00
65	c	n-Nitrosodiphenylamine	0.593	0.622	-4.9	71	-0.01
66		4-Bromophenyl-phenylether	0.208	0.197	5.3	64	-0.01
67		Hexachlorobenzene	0.222	0.234	-5.4	69	0.00
68		Atrazine	0.191	0.164	14.1	58	-0.01
69	C	Pentachlorophenol	0.109	0.116	-6.4	63	0.00
70		Phenanthrene	1.084	1.138	-5.0	68	0.00
71		Anthracene	1.086	1.026	5.5	71	-0.01
72		Carbazole	1.005	0.980	2.5	72	-0.01
73		Di-n-butylphthalate	1.174	1.233	-5.0	76	0.00
74	C	Fluoranthene	1.082	1.065	1.6	74	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	70	0.00
76		Benzidine	0.580	0.573	1.2	73	-0.01
77		Pyrene	1.467	1.529	-4.2	79	0.00
78	S	Terphenyl-d14	0.825	0.817	1.0	77	-0.01
79		Butylbenzylphthalate	0.667	0.650	2.5	64	-0.01
80		Benzo(a)anthracene	1.129	1.207	-6.9	76	0.00
81		3,3'-Dichlorobenzidine	0.428	0.451	-5.4	79	-0.01
82		Chrysene	1.132	1.143	-1.0	79	0.00
83		Bis(2-ethylhexyl)phthalate	0.786	0.828	-5.3	75	0.00
84	c	Di-n-octyl phthalate	1.456	1.550	-6.5	76	0.00
85		Indeno(1,2,3-cd)pyrene	0.981	1.056	-7.6	78	0.00
86	I	Perylene-d12	1.000	1.000	0.0	79	0.00
87		Benzo(b)fluoranthene	1.245	1.368	-9.9	82	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF010917\
Data File : BF092141.D
Acq On : 9 Jan 2017 17:00
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 09 22:40:52 2017
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 06 15:43:51 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	1.062	0.935	12.0	72	0.00
89 C	Benzo(a)pyrene	1.105	1.154	-4.4	81	0.00
90	Dibenzo(a,h)anthracene	0.908	0.917	-1.0	78	0.00
91	Benzo(g,h,i)perylene	0.945	0.926	2.0	76	0.00

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