Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100416\
Data File : BF090369.D
Acq On : 5 Oct 2016 3:45
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 05 04:50:36 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF100416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 05 03:16:32 2016
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	65	0.00
2		1,4-Dioxane	0.619	0.774	-25.0#	80	0.00
3		Pyridine	1.641	2.035	-24.0	79	0.02
4		n-Nitrosodimethylamine	0.594	0.798	-34.3#	85	0.00
5	S	2-Fluorophenol	1.268	1.535	-21.1	72	0.00
6		Aniline	2.178	2.318	-6.4	67	0.00
7	S	Phenol-d6	1.585	1.757	-10.9	68	0.00
8		2-Chlorophenol	1.425	1.487	-4.4	69	0.00
9		Benzaldehyde	0.736	0.855	-16.2	76	0.00
10	C	Phenol	1.928	2.147	-11.4	65	0.00
11		bis(2-Chloroethyl)ether	1.369	1.373	-0.3	59	0.00
12		1,3-Dichlorobenzene	1.480	1.548	-4.6	70	0.00
13	C	1,4-Dichlorobenzene	1.503	1.457	3.1	61	0.00
14		1,2-Dichlorobenzene	1.429	1.336	6.5	61	0.00
15		Benzyl Alcohol	1.117	1.329	-19.0	79	0.00
16		2,2'-oxybis(1-Chloropropane	2.092	2.484	-18.7	74	0.00
17		2-Methylphenol	1.104	1.188	-7.6	66	0.00
18		Hexachloroethane	0.544	0.363	33.3#	41#	0.00
19	P	n-Nitroso-di-n-propylamine	1.012	1.247	-23.2	83	0.00
20		3+4-Methylphenols	1.403	1.454	-3.6	68	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	63	0.00
22		Acetophenone	0.463	0.513	-10.8	71	0.00
23	S	Nitrobenzene-d5	0.343	0.375	-9.3	63	0.00
24		Nitrobenzene	0.366	0.452	-23.5	82	0.00
25		Isophorone	0.653	0.761	-16.5	77	0.00
26	C	2-Nitrophenol	0.141	0.161	-14.2	71	0.00
27		2,4-Dimethylphenol	0.319	0.331	-3.8	63	0.00
28		bis(2-Chloroethoxy)methane	0.405	0.413	-2.0	59	0.00
29	C	2,4-Dichlorophenol	0.261	0.260	0.4	59	0.00
30		1,2,4-Trichlorobenzene	0.292	0.254	13.0	54	0.00
31		Naphthalene	0.956	0.968	-1.3	67	0.00
32		Benzoic acid	0.188	0.234	-24.5	75	0.00
33		4-Chloroaniline	0.382	0.387	-1.3	58	0.00
34	C	Hexachlorobutadiene	0.172	0.135	21.5#	49#	0.00
35		Caprolactam	0.068	0.079	-16.2	69	0.00
36	C	4-Chloro-3-methylphenol	0.280	0.298	-6.4	69	0.00
37		2-Methylnaphthalene	0.600	0.539	10.2	54	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	60	0.00
39		1,2,4,5-Tetrachlorobenzene	0.565	0.535	5.3	59	0.00
40	P	Hexachlorocyclopentadiene	0.330	0.054	83.6#	10#	0.00
41	S	2,4,6-Tribromophenol	0.174	0.180	-3.4	60	0.00
42	C	2,4,6-Trichlorophenol	0.355	0.326	8.2	56	0.00
43		2,4,5-Trichlorophenol	0.357	0.345	3.4	56	0.00
44	S	2-Fluorobiphenyl	1.227	1.290	-5.1	63	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100416\
Data File : BF090369.D
Acq On : 5 Oct 2016 3:45
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 05 04:50:36 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF100416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 05 03:16:32 2016
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	1.548	1.688	-9.0	72	0.00
46		2-Chloronaphthalene	1.162	1.124	3.3	59	0.00
47		2-Nitroaniline	0.317	0.490	-54.6#	93	0.00
48		Acenaphthylene	1.758	1.830	-4.1	60	0.00
49		Dimethylphthalate	1.184	1.353	-14.3	72	0.00
50		2,6-Dinitrotoluene	0.259	0.275	-6.2	72	0.00
51	C	Acenaphthene	1.070	1.103	-3.1	63	0.00
52		3-Nitroaniline	0.293	0.377	-28.7#	77	0.00
53	P	2,4-Dinitrophenol	0.078	0.032#	59.0#	22#	0.00
54		Dibenzofuran	1.415	1.505	-6.4	65	0.00
55	P	4-Nitrophenol	0.234	0.316	-35.0#	77	0.00
56		2,4-Dinitrotoluene	0.312	0.338	-8.3	71	0.00
57		Fluorene	1.183	1.344	-13.6	70	0.00
58		2,3,4,6-Tetrachlorophenol	0.252	0.266	-5.6	59	0.00
59		Diethylphthalate	1.210	1.299	-7.4	64	0.00
60		4-Chlorophenyl-phenylether	0.561	0.567	-1.1	62	0.00
61		4-Nitroaniline	0.264	0.353	-33.7#	76	0.00
62		Azobenzene	1.261	1.588	-25.9#	84	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	73	0.00
64		4,6-Dinitro-2-methylphenol	0.076	0.036	52.6#	31#	0.00
65	c	n-Nitrosodiphenylamine	0.682	0.619	9.2	64	0.00
66		4-Bromophenyl-phenylether	0.216	0.190	12.0	66	0.00
67		Hexachlorobenzene	0.254	0.210	17.3	63	0.00
68		Atrazine	0.163	0.146	10.4	69	0.00
69	C	Pentachlorophenol	0.142	0.132	7.0	65	0.00
70		Phenanthrene	1.073	0.998	7.0	67	0.00
71		Anthracene	1.071	1.144	-6.8	75	0.00
72		Carbazole	0.985	1.059	-7.5	81	0.00
73		Di-n-butylphthalate	1.177	1.277	-8.5	75	0.00
74	C	Fluoranthene	1.018	1.124	-10.4	76	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	94	0.00
76		Benzidine	0.550	0.583	-6.0	94	0.00
77		Pyrene	1.281	1.171	8.6	82	0.00
78	S	Terphenyl-d14	0.859	0.698	18.7	74	0.00
79		Butylbenzylphthalate	0.505	0.634	-25.5#	113	0.00
80		Benzo(a)anthracene	1.121	1.121	0.0	97	0.00
81		3,3'-Dichlorobenzidine	0.386	0.427	-10.6	103	0.00
82		Chrysene	1.051	0.933	11.2	91	0.00
83		Bis(2-ethylhexyl)phthalate	0.807	0.913	-13.1	105	0.00
84	c	Di-n-octyl phthalate	1.173	1.536	-30.9#	118	0.00
85		Indeno(1,2,3-cd)pyrene	1.129	0.769	31.9#	68	0.00
86	I	Perylene-d12	1.000	1.000	0.0	87	0.00
87		Benzo(b)fluoranthene	1.150	1.273	-10.7	97	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF100416\
Data File : BF090369.D
Acq On : 5 Oct 2016 3:45
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 05 04:50:36 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF100416.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 05 03:16:32 2016
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	1.099	1.289	-17.3	92	0.00
89 C	Benzo(a)pyrene	1.049	1.108	-5.6	87	0.00
90	Dibenzo(a,h)anthracene	1.029	0.858	16.6	69	0.02
91	Benzo(g,h,i)perylene	1.008	0.778	22.8	64	0.00

(#) = Out of Range SPCC's out = 1 CCC's out = 2