Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_F\DATA\BF031016\
Data File : BF085354.D
Acq On : 11 Mar 2016 1:16
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 11 02:47:05 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 11 00:01:03 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	61	-0.05
2		1,4-Dioxane	0.610	0.643	-5.4	67	-0.06
3		Pyridine	1.526	1.450	5.0	60	-0.07
4		n-Nitrosodimethylamine	0.653	0.639	2.1	59	-0.07
5	S	2-Fluorophenol	1.192	1.323	-11.0	75	-0.03
6		Aniline	2.189	2.078	5.1	58	-0.03
7	S	Phenol-d6	1.562	1.539	1.5	64	-0.03
8		2-Chlorophenol	1.435	1.391	3.1	61	-0.05
9		Benzaldehyde	0.804	0.753	6.3	65	-0.03
10	C	Phenol	1.673	1.860	-11.2	72	-0.03
11		bis(2-Chloroethyl)ether	1.390	1.341	3.5	56	-0.05
12		1,3-Dichlorobenzene	1.541	1.537	0.3	62	-0.05
13	C	1,4-Dichlorobenzene	1.545	1.652	-6.9	71	-0.03
14		1,2-Dichlorobenzene	1.401	1.313	6.3	56	-0.05
15		Benzyl Alcohol	1.147	1.013	11.7	56	-0.05
16		2,2'-oxybis(1-Chloropropane	2.003	1.792	10.5	58	-0.05
17		2-Methylphenol	1.171	1.072	8.5	55	-0.05
18		Hexachloroethane	0.570	0.527	7.5	57	-0.05
19	P	n-Nitroso-di-n-propylamine	1.018	0.899	11.7	54	-0.05
20		3+4-Methylphenols	1.387	1.393	-0.4	67	-0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	62	-0.03
22		Acetophenone	0.488	0.434	11.1	55	-0.05
23	S	Nitrobenzene-d5	0.374	0.346	7.5	56	-0.03
24		Nitrobenzene	0.380	0.325	14.5	52	-0.05
25		Isophorone	0.693	0.611	11.8	54	-0.05
26	C	2-Nitrophenol	0.185	0.188	-1.6	60	-0.05
27		2,4-Dimethylphenol	0.338	0.317	6.2	55	-0.05
28		bis(2-Chloroethoxy)methane	0.386	0.389	-0.8	63	-0.03
29	C	2,4-Dichlorophenol	0.272	0.265	2.6	59	-0.05
30		1,2,4-Trichlorobenzene	0.294	0.293	0.3	62	-0.03
31		Naphthalene	0.983	0.974	0.9	66	-0.03
32		Benzoic acid	0.224	0.196	12.5	49#	-0.07
33		4-Chloroaniline	0.398	0.395	0.8	65	-0.03
34	C	Hexachlorobutadiene	0.153	0.160	-4.6	63	-0.05
35		Caprolactam	0.089	0.071	20.2	49#	-0.06
36	C	4-Chloro-3-methylphenol	0.309	0.245	20.7#	50#	-0.05
37		2-Methylnaphthalene	0.631	0.537	14.9	52	-0.05
38	I	Acenaphthene-d10	1.000	1.000	0.0	41#	-0.05
39		1,2,4,5-Tetrachlorobenzene	0.572	0.711	-24.3	56	-0.03
40	P	Hexachlorocyclopentadiene	0.308	0.215	30.2#	28#	-0.03
41	S	2,4,6-Tribromophenol	0.171	0.177	-3.5	41#	-0.05
42	C	2,4,6-Trichlorophenol	0.426	0.451	-5.9	44#	-0.05
43		2,4,5-Trichlorophenol	0.404	0.443	-9.7	44#	-0.03
44	S	2-Fluorobiphenyl	1.315	1.523	-15.8	51	-0.05

Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_F\DATA\BF031016\
Data File : BF085354.D
Acq On : 11 Mar 2016 1:16
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 11 02:47:05 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 11 00:01:03 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	1.709	1.967	-15.1	53	-0.03
46		2-Chloronaphthalene	1.302	1.459	-12.1	47#	-0.05
47		2-Nitroaniline	0.404	0.420	-4.0	42#	-0.03
48		Acenaphthylene	2.027	2.063	-1.8	44#	-0.05
49		Dimethylphthalate	1.535	1.696	-10.5	45#	-0.05
50		2,6-Dinitrotoluene	0.328	0.337	-2.7	41#	-0.05
51	C	Acenaphthene	1.231	1.207	1.9	43#	-0.05
52		3-Nitroaniline	0.393	0.416	-5.9	40#	-0.05
53	P	2,4-Dinitrophenol	0.137	0.095	30.7#	27#	-0.03
54		Dibenzofuran	1.613	1.598	0.9	42#	-0.05
55	P	4-Nitrophenol	0.309	0.263	14.9	32#	-0.05
56		2,4-Dinitrotoluene	0.420	0.371	11.7	36#	-0.05
57		Fluorene	1.267	1.269	-0.2	41#	-0.05
58		2,3,4,6-Tetrachlorophenol	0.284	0.279	1.8	39#	-0.05
59		Diethylphthalate	1.490	1.560	-4.7	44#	-0.05
60		4-Chlorophenyl-phenylether	0.616	0.624	-1.3	44#	-0.05
61		4-Nitroaniline	0.367	0.304	17.2	33#	-0.05
62		Azobenzene	1.468	1.166	20.6	32#	-0.05
63	I	Phenanthrene-d10	1.000	1.000	0.0	36#	-0.05
64		4,6-Dinitro-2-methylphenol	0.120	0.094	21.7	25#	-0.04
65	c	n-Nitrosodiphenylamine	0.622	0.713	-14.6	44#	-0.03
66		4-Bromophenyl-phenylether	0.194	0.223	-14.9	41#	-0.05
67		Hexachlorobenzene	0.207	0.225	-8.7	39#	-0.05
68		Atrazine	0.173	0.201	-16.2	49#	-0.03
69	C	Pentachlorophenol	0.115	0.135	-17.4	42#	-0.03
70		Phenanthrene	1.048	1.044	0.4	40#	-0.05
71		Anthracene	1.113	1.168	-4.9	41#	-0.05
72		Carbazole	0.976	0.923	5.4	35#	-0.05
73		Di-n-butylphthalate	1.233	1.272	-3.2	40#	-0.03
74	C	Fluoranthene	1.052	1.057	-0.5	38#	-0.05
75	I	Chrysene-d12	1.000	1.000	0.0	58	-0.05
76		Benzidine	0.595	0.589	1.0	49#	-0.03
77		Pyrene	1.505	1.126	25.2#	39#	-0.05
78	S	Terphenyl-d14	0.862	0.701	18.7	46#	-0.03
79		Butylbenzylphthalate	0.683	0.535	21.7	42#	-0.03
80		Benzo(a)anthracene	1.197	1.102	7.9	52	-0.03
81		3,3'-Dichlorobenzidine	0.378	0.435	-15.1	62	-0.03
82		Chrysene	1.106	0.953	13.8	45#	-0.05
83		Bis(2-ethylhexyl)phthalate	0.899	0.749	16.7	46#	-0.03
84	c	Di-n-octyl phthalate	1.369	1.406	-2.7	53	-0.03
85		Indeno(1,2,3-cd)pyrene	0.863	1.273	-47.5#	72	-0.07
86	I	Perylene-d12	1.000	1.000	0.0	72	-0.05
87		Benzo(b)fluoranthene	1.333	1.208	9.4	62	-0.03

Evaluate Continuing Calibration Report

Data Path : S:\HPCHEM1\BNA_F\DATA\BF031016\
Data File : BF085354.D
Acq On : 11 Mar 2016 1:16
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Mar 11 02:47:05 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF030316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 11 00:01:03 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	1.075	1.043	3.0	81	-0.05
89 C	Benzo(a)pyrene	1.105	1.111	-0.5	72	-0.05
90	Dibenzo(a,h)anthracene	0.943	1.055	-11.9	73	-0.06
91	Benzo(g,h,i)perylene	0.948	0.990	-4.4	67	-0.07

(#) = Out of Range SPCC's out = 0 CCC's out = 1