Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF041817\
Data File : BF094488.D
Acq On : 18 Apr 2017 15:55
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 19 01:52:14 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Apr 17 14:59:23 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	77	0.00
2		1,4-Dioxane	0.701	0.778	-11.0	86	0.02
3		Pyridine	1.532	1.676	-9.4	80	0.03
4		n-Nitrosodimethylamine	0.634	0.648	-2.2	78	0.01
5	S	2-Fluorophenol	1.205	1.340	-11.2	87	0.01
6		Aniline	1.889	1.887	0.1	79	0.01
7	S	Phenol-d6	1.421	1.435	-1.0	80	0.00
8		2-Chlorophenol	1.372	1.414	-3.1	80	0.00
9		Benzaldehyde	1.081	1.062	1.8	77	0.01
10	C	Phenol	1.746	1.751	-0.3	80	0.00
11		bis(2-Chloroethyl)ether	1.275	1.309	-2.7	80	0.00
12		1,3-Dichlorobenzene	1.520	1.562	-2.8	81	0.01
13	C	1,4-Dichlorobenzene	1.529	1.585	-3.7	81	0.00
14		1,2-Dichlorobenzene	1.408	1.477	-4.9	83	0.00
15		Benzyl Alcohol	1.054	1.088	-3.2	80	0.00
16		2,2'-oxybis(1-Chloropropane	1.672	1.770	-5.9	83	0.00
17		2-Methylphenol	1.080	1.115	-3.2	81	0.00
18		Hexachloroethane	0.524	0.554	-5.7	82	0.00
19	P	n-Nitroso-di-n-propylamine	0.942	0.994	-5.5	82	0.00
20		3+4-Methylphenols	1.468	1.535	-4.6	81	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	86	0.00
22		Acetophenone	0.486	0.449	7.6	80	0.00
23	S	Nitrobenzene-d5	0.324	0.304	6.2	81	0.00
24		Nitrobenzene	0.341	0.322	5.6	80	0.00
25		Isophorone	0.600	0.581	3.2	83	0.00
26	C	2-Nitrophenol	0.183	0.178	2.7	82	0.00
27		2,4-Dimethylphenol	0.298	0.291	2.3	84	0.00
28		bis(2-Chloroethoxy)methane	0.388	0.381	1.8	85	0.00
29	C	2,4-Dichlorophenol	0.277	0.285	-2.9	87	0.00
30		1,2,4-Trichlorobenzene	0.286	0.283	1.0	86	0.00
31		Naphthalene	0.961	0.958	0.3	87	0.00
32		Benzoic acid	0.157	0.167	-6.4	91	0.00
33		4-Chloroaniline	0.400	0.407	-1.7	87	0.00
34	C	Hexachlorobutadiene	0.143	0.145	-1.4	87	0.00
35		Caprolactam	0.087	0.094	-8.0	93	0.00
36	C	4-Chloro-3-methylphenol	0.290	0.276	4.8	82	0.00
37		2-Methylnaphthalene	0.658	0.664	-0.9	87	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	86	0.00
39		1,2,4,5-Tetrachlorobenzene	0.570	0.567	0.5	88	0.00
40	P	Hexachlorocyclopentadiene	0.230	0.257	-11.7	93	0.00
41	S	2,4,6-Tribromophenol	0.156	0.168	-7.7	93	0.00
42	C	2,4,6-Trichlorophenol	0.400	0.389	2.8	84	0.00
43		2,4,5-Trichlorophenol	0.411	0.397	3.4	83	0.00
44	S	2-Fluorobiphenyl	1.359	1.270	6.5	84	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF041817\
Data File : BF094488.D
Acq On : 18 Apr 2017 15:55
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 19 01:52:14 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Apr 17 14:59:23 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	1.634	1.633	0.1	87	0.00
46		2-Chloronaphthalene	1.299	1.306	-0.5	89	0.00
47		2-Nitroaniline	0.374	0.387	-3.5	88	0.00
48		Acenaphthylene	2.025	1.973	2.6	86	0.00
49		Dimethylphthalate	1.490	1.552	-4.2	91	0.00
50		2,6-Dinitrotoluene	0.335	0.345	-3.0	88	0.00
51	C	Acenaphthene	1.213	1.189	2.0	87	0.00
52		3-Nitroaniline	0.387	0.393	-1.6	88	0.00
53	P	2,4-Dinitrophenol	0.158	0.171	-8.2	91	0.00
54		Dibenzofuran	1.763	1.788	-1.4	91	0.00
55	P	4-Nitrophenol	0.273	0.274	-0.4	92	0.00
56		2,4-Dinitrotoluene	0.410	0.443	-8.0	94	0.00
57		Fluorene	1.373	1.411	-2.8	91	0.00
58		2,3,4,6-Tetrachlorophenol	0.294	0.317	-7.8	92	0.00
59		Diethylphthalate	1.406	1.513	-7.6	95	0.00
60		4-Chlorophenyl-phenylether	0.571	0.616	-7.9	94	0.00
61		4-Nitroaniline	0.393	0.369	6.1	80	0.00
62		Azobenzene	1.365	1.444	-5.8	92	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
64		4,6-Dinitro-2-methylphenol	0.131	0.145	-10.7	93	0.00
65	c	n-Nitrosodiphenylamine	0.696	0.729	-4.7	92	0.00
66		4-Bromophenyl-phenylether	0.199	0.212	-6.5	93	0.00
67		Hexachlorobenzene	0.200	0.202	-1.0	89	0.00
68		Atrazine	0.208	0.221	-6.3	93	0.00
69	C	Pentachlorophenol	0.079	0.092	-16.5	98	0.00
70		Phenanthrene	1.126	1.129	-0.3	90	0.00
71		Anthracene	1.165	1.173	-0.7	88	0.00
72		Carbazole	1.093	1.115	-2.0	90	0.00
73		Di-n-butylphthalate	1.204	1.329	-10.4	96	0.00
74	C	Fluoranthene	1.167	1.219	-4.5	92	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	99	0.00
76		Benzidine	0.810	0.739	8.8	87	0.00
77		Pyrene	1.397	1.303	6.7	92	0.00
78	S	Terphenyl-d14	0.748	0.705	5.7	96	0.00
79		Butylbenzylphthalate	0.612	0.632	-3.3	101	0.00
80		Benzo(a)anthracene	1.162	1.175	-1.1	100	0.00
81		3,3'-Dichlorobenzidine	0.412	0.424	-2.9	99	0.00
82		Chrysene	1.062	1.091	-2.7	104	0.00
83		Bis(2-ethylhexyl)phthalate	0.822	0.938	-14.1	112	0.00
84	c	Di-n-octyl phthalate	1.379	1.546	-12.1	109	0.00
85		Indeno(1,2,3-cd)pyrene	1.011	0.969	4.2	99	0.00
86	I	Perylene-d12	1.000	1.000	0.0	99	0.00
87		Benzo(b)fluoranthene	1.223	1.253	-2.5	100	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF041817\
Data File : BF094488.D
Acq On : 18 Apr 2017 15:55
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 19 01:52:14 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Apr 17 14:59:23 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	1.158	1.169	-0.9	103	0.00
89 C	Benzo(a)pyrene	1.138	1.144	-0.5	100	0.00
90	Dibenzo(a,h)anthracene	0.945	0.914	3.3	100	0.00
91	Benzo(g,h,i)perylene	0.962	0.867	9.9	95	0.00

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