Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091815\
Data File : BF081694.D
Acq On : 18 Sep 2015 12:56
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 18 23:19:59 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 15 14:03:59 2015
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	118	-0.04
2		1,4-Dioxane	0.579	1.356	-134.2#	280#	-0.03
3		Pyridine	1.596	1.477	7.5	95	-0.03
4		n-Nitrosodimethylamine	0.887	1.052	-18.6	132	-0.03
5	S	2-Fluorophenol	1.289	1.269	1.6	121	-0.03
6		Aniline	2.410	2.380	1.2	117	-0.03
7	S	Phenol-d6	1.706	1.752	-2.7	119	-0.03
8		2-Chlorophenol	1.420	1.450	-2.1	121	-0.04
9		Benzaldehyde	1.069	0.922	13.8	101	-0.03
10	C	Phenol	1.979	1.861	6.0	101	-0.03
11		bis(2-Chloroethyl)ether	1.428	1.457	-2.0	120	-0.04
12		1,3-Dichlorobenzene	1.518	1.499	1.3	119	-0.04
13	C	1,4-Dichlorobenzene	1.545	1.549	-0.3	119	-0.04
14		1,2-Dichlorobenzene	1.391	1.451	-4.3	125	-0.03
15		Benzyl Alcohol	1.400	1.477	-5.5	117	-0.03
16		2,2'-oxybis(1-Chloropropane	2.736	3.059	-11.8	138	-0.04
17		2-Methylphenol	1.133	1.179	-4.1	121	-0.03
18		Hexachloroethane	0.601	0.641	-6.7	126	-0.04
19	P	n-Nitroso-di-n-propylamine	1.161	1.295	-11.5	138	-0.02
20		3+4-Methylphenols	1.528	1.560	-2.1	122	-0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	128	-0.04
22		Acetophenone	0.546	0.537	1.6	122	-0.03
23	S	Nitrobenzene-d5	0.415	0.432	-4.1	130	-0.03
24		Nitrobenzene	0.430	0.457	-6.3	129	-0.03
25		Isophorone	0.771	0.805	-4.4	135	-0.03
26	C	2-Nitrophenol	0.188	0.185	1.6	132	-0.04
27		2,4-Dimethylphenol	0.324	0.318	1.9	112	-0.04
28		bis(2-Chloroethoxy)methane	0.445	0.454	-2.0	116	-0.03
29	C	2,4-Dichlorophenol	0.281	0.285	-1.4	125	-0.03
30		1,2,4-Trichlorobenzene	0.305	0.316	-3.6	126	-0.04
31		Naphthalene	1.067	1.065	0.2	127	-0.03
32		Benzoic acid	0.221	0.172	22.2	90	0.00
33		4-Chloroaniline	0.435	0.448	-3.0	115	-0.03
34	C	Hexachlorobutadiene	0.186	0.204	-9.7	145	-0.04
35		Caprolactam	0.086	0.086	0.0	122	-0.02
36	C	4-Chloro-3-methylphenol	0.315	0.353	-12.1	143	-0.03
37		2-Methylnaphthalene	0.694	0.714	-2.9	141	-0.04
38	I	Acenaphthene-d10	1.000	1.000	0.0	132	-0.04
39		1,2,4,5-Tetrachlorobenzene	0.633	0.634	-0.2	129	-0.04
40	P	Hexachlorocyclopentadiene	0.310	0.257	17.1	107	-0.06
41	S	2,4,6-Tribromophenol	0.178	0.188	-5.6	133	-0.04
42	C	2,4,6-Trichlorophenol	0.437	0.442	-1.1	136	-0.04
43		2,4,5-Trichlorophenol	0.445	0.443	0.4	128	-0.04
44	S	2-Fluorobiphenyl	1.561	1.603	-2.7	130	-0.04

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091815\
Data File : BF081694.D
Acq On : 18 Sep 2015 12:56
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 18 23:19:59 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 15 14:03:59 2015
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	1.840	1.783	3.1	142	-0.04
46		2-Chloronaphthalene	1.329	1.326	0.2	123	-0.04
47		2-Nitroaniline	0.542	0.582	-7.4	137	-0.04
48		Acenaphthylene	2.357	2.335	0.9	143	-0.06
49		Dimethylphthalate	1.554	1.621	-4.3	143	-0.04
50		2,6-Dinitrotoluene	0.353	0.342	3.1	121	-0.04
51	C	Acenaphthene	1.341	1.292	3.7	142	-0.04
52		3-Nitroaniline	0.402	0.388	3.5	125	-0.04
53	P	2,4-Dinitrophenol	0.149	0.146	2.0	122	-0.03
54		Dibenzofuran	1.958	1.959	-0.1	144	-0.04
55	P	4-Nitrophenol	0.252	0.228	9.5	102	-0.03
56		2,4-Dinitrotoluene	0.449	0.463	-3.1	135	-0.03
57		Fluorene	1.486	1.586	-6.7	133	-0.04
58		2,3,4,6-Tetrachlorophenol	0.340	0.345	-1.5	139	-0.04
59		Diethylphthalate	1.635	1.782	-9.0	143	-0.04
60		4-Chlorophenyl-phenylether	0.693	0.760	-9.7	151#	-0.04
61		4-Nitroaniline	0.406	0.369	9.1	128	-0.03
62		Azobenzene	1.817	1.931	-6.3	132	-0.04
63	I	Phenanthrene-d10	1.000	1.000	0.0	139	-0.04
64		4,6-Dinitro-2-methylphenol	0.112	0.110	1.8	118	-0.04
65	c	n-Nitrosodiphenylamine	0.728	0.681	6.5	130	-0.04
66		4-Bromophenyl-phenylether	0.202	0.209	-3.5	161#	-0.04
67		Hexachlorobenzene	0.211	0.211	0.0	148	-0.06
68		Atrazine	0.211	0.201	4.7	136	-0.04
69	C	Pentachlorophenol	0.082	0.081	1.2	136	-0.04
70		Phenanthrene	1.112	1.110	0.2	134	-0.06
71		Anthracene	1.142	1.100	3.7	138	-0.06
72		Carbazole	1.072	1.028	4.1	142	-0.04
73		Di-n-butylphthalate	1.266	1.418	-12.0	165#	-0.06
74	C	Fluoranthene	1.204	1.181	1.9	146	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	120	-0.02
76		Benzidine	0.859	0.730	15.0	116	-0.02
77		Pyrene	1.481	1.557	-5.1	130	-0.04
78	S	Terphenyl-d14	0.859	0.981	-14.2	165#	-0.03
79		Butylbenzylphthalate	0.644	0.747	-16.0	152#	-0.03
80		Benzo(a)anthracene	1.274	1.269	0.4	123	-0.03
81		3,3'-Dichlorobenzidine	0.430	0.407	5.3	130	-0.03
82		Chrysene	1.142	1.105	3.2	147	-0.03
83		Bis(2-ethylhexyl)phthalate	0.850	1.040	-22.4	169#	-0.03
84	c	Di-n-octyl phthalate	1.400	1.500	-7.1	144	-0.03
85		Indeno(1,2,3-cd)pyrene	0.838	0.839	-0.1	132	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	123	-0.02
87		Benzo(b)fluoranthene	1.428	1.241	13.1	87	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF091815\
Data File : BF081694.D
Acq On : 18 Sep 2015 12:56
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 18 23:19:59 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Sep 15 14:03:59 2015
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	1.185	1.350	-13.9	169#	-0.02
89 C	Benzo(a)pyrene	1.210	1.213	-0.2	131	-0.02
90	Dibenzo(a,h)anthracene	0.935	1.016	-8.7	133	-0.04
91	Benzo(g,h,i)perylene	0.892	0.929	-4.1	131	-0.04

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