Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF081815\
Data File : BF081060.D
Acq On : 19 Aug 2015 00:36
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 19 01:56:18 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 17 17:57:25 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	101	-0.02
2		1,4-Dioxane	0.627	0.150	76.1#	24#	-0.06
3		Pyridine	1.769	1.808	-2.2	92	-0.09
4		n-Nitrosodimethylamine	0.985	0.943	4.3	95	-0.05
5	S	2-Fluorophenol	1.330	1.229	7.6	97	-0.02
6		Aniline	2.523	2.618	-3.8	107	-0.01
7	S	Phenol-d6	1.735	1.711	1.4	94	-0.01
8		2-Chlorophenol	1.455	1.557	-7.0	100	-0.01
9		Benzaldehyde	1.118	1.198	-7.2	101	-0.01
10	C	Phenol	2.049	1.848	9.8	82	-0.01
11		bis(2-Chloroethyl)ether	1.572	1.722	-9.5	109	-0.01
12		1,3-Dichlorobenzene	1.564	1.669	-6.7	107	-0.01
13	C	1,4-Dichlorobenzene	1.549	1.700	-9.7	112	-0.01
14		1,2-Dichlorobenzene	1.449	1.453	-0.3	94	-0.02
15		Benzyl Alcohol	1.404	1.548	-10.3	120	-0.01
16		2,2'-oxybis(1-Chloropropane	3.089	3.151	-2.0	99	-0.01
17		2-Methylphenol	1.249	1.222	2.2	95	-0.01
18		Hexachloroethane	0.626	0.657	-5.0	102	-0.02
19	P	n-Nitroso-di-n-propylamine	1.202	1.235	-2.7	97	-0.01
20		3+4-Methylphenols	1.585	1.582	0.2	102	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	109	-0.01
22		Acetophenone	0.525	0.495	5.7	96	-0.01
23	S	Nitrobenzene-d5	0.438	0.421	3.9	105	-0.01
24		Nitrobenzene	0.458	0.464	-1.3	105	-0.01
25		Isophorone	0.816	0.837	-2.6	110	-0.01
26	C	2-Nitrophenol	0.190	0.208	-9.5	124	-0.01
27		2,4-Dimethylphenol	0.337	0.318	5.6	93	-0.01
28		bis(2-Chloroethoxy)methane	0.482	0.482	0.0	109	-0.01
29	C	2,4-Dichlorophenol	0.284	0.304	-7.0	110	-0.01
30		1,2,4-Trichlorobenzene	0.315	0.312	1.0	102	-0.02
31		Naphthalene	1.115	1.196	-7.3	115	-0.01
32		Benzoic acid	0.189	0.206	-9.0	138	0.05
33		4-Chloroaniline	0.470	0.500	-6.4	124	-0.01
34	C	Hexachlorobutadiene	0.178	0.189	-6.2	114	-0.01
35		Caprolactam	0.099	0.096	3.0	101	0.02
36	C	4-Chloro-3-methylphenol	0.338	0.348	-3.0	110	0.00
37		2-Methylnaphthalene	0.704	0.704	0.0	105	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	115	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.645	0.660	-2.3	117	-0.01
40	P	Hexachlorocyclopentadiene	0.334	0.353	-5.7	113	-0.01
41	S	2,4,6-Tribromophenol	0.173	0.173	0.0	119	-0.02
42	C	2,4,6-Trichlorophenol	0.456	0.485	-6.4	122	-0.01
43		2,4,5-Trichlorophenol	0.456	0.477	-4.6	107	-0.01
44	S	2-Fluorobiphenyl	1.526	1.335	12.5	111	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF081815\
Data File : BF081060.D
Acq On : 19 Aug 2015 00:36
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 19 01:56:18 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 17 17:57:25 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	1.762	1.918	-8.9	113	-0.01
46		2-Chloronaphthalene	1.415	1.370	3.2	114	-0.01
47		2-Nitroaniline	0.579	0.524	9.5	102	-0.01
48		Acenaphthylene	2.532	2.329	8.0	112	-0.02
49		Dimethylphthalate	1.647	1.682	-2.1	109	-0.01
50		2,6-Dinitrotoluene	0.354	0.377	-6.5	112	-0.01
51	C	Acenaphthene	1.516	1.483	2.2	114	-0.01
52		3-Nitroaniline	0.443	0.397	10.4	105	-0.01
53	P	2,4-Dinitrophenol	0.167	0.204	-22.2	134	-0.02
54		Dibenzofuran	2.031	1.921	5.4	115	-0.02
55	P	4-Nitrophenol	0.311	0.354	-13.8	137	-0.01
56		2,4-Dinitrotoluene	0.469	0.540	-15.1	128	-0.01
57		Fluorene	1.562	1.471	5.8	114	-0.02
58		2,3,4,6-Tetrachlorophenol	0.353	0.397	-12.5	117	-0.02
59		Diethylphthalate	1.743	1.958	-12.3	125	-0.01
60		4-Chlorophenyl-phenylether	0.661	0.713	-7.9	121	-0.02
61		4-Nitroaniline	0.452	0.518	-14.6	134	0.00
62		Azobenzene	2.009	2.232	-11.1	118	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	120	-0.02
64		4,6-Dinitro-2-methylphenol	0.119	0.126	-5.9	114	-0.01
65	c	n-Nitrosodiphenylamine	0.714	0.758	-6.2	123	-0.01
66		4-Bromophenyl-phenylether	0.196	0.194	1.0	122	-0.02
67		Hexachlorobenzene	0.199	0.220	-10.6	124	-0.01
68		Atrazine	0.199	0.192	3.5	116	-0.01
69	C	Pentachlorophenol	0.119	0.136	-14.3	135	-0.01
70		Phenanthrene	1.185	1.081	8.8	112	-0.02
71		Anthracene	1.153	1.241	-7.6	125	-0.01
72		Carbazole	1.110	1.259	-13.4	121	-0.01
73		Di-n-butylphthalate	1.344	1.591	-18.4	137	-0.02
74	C	Fluoranthene	1.279	1.207	5.6	109	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	119	-0.01
76		Benzidine	0.756	1.052	-39.2#	134	-0.02
77		Pyrene	1.626	1.688	-3.8	130	-0.02
78	S	Terphenyl-d14	0.933	1.104	-18.3	134	-0.01
79		Butylbenzylphthalate	0.699	0.978	-39.9#	150	-0.02
80		Benzo(a)anthracene	1.341	1.594	-18.9	133	-0.02
81		3,3'-Dichlorobenzidine	0.434	0.491	-13.1	134	-0.02
82		Chrysene	1.209	1.485	-22.8	135	-0.01
83		Bis(2-ethylhexyl)phthalate	0.895	1.299	-45.1#	164#	-0.02
84	c	Di-n-octyl phthalate	1.361	2.131	-56.6#	172#	-0.02
85		Indeno(1,2,3-cd)pyrene	0.849	1.205	-41.9#	160#	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	160#	-0.02
87		Benzo(b)fluoranthene	1.415	1.656	-17.0	170#	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF081815\
Data File : BF081060.D
Acq On : 19 Aug 2015 00:36
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 19 01:56:18 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF081715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 17 17:57:25 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	1.318	1.002	24.0	132	-0.02
89 C	Benzo(a)pyrene	1.233	1.257	-1.9	161#	-0.01
90	Dibenzo(a,h)anthracene	0.960	0.996	-3.8	161#	-0.03
91	Benzo(g,h,i)perylene	0.974	0.969	0.5	156#	-0.02

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