Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122015\
Data File : BF083949.D
Acq On : 19 Dec 2015 12:42
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 21 02:59:21 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 18 14:23:22 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	111	-0.02
2		1,4-Dioxane	0.554	0.470	15.2	95	-0.05
3		Pyridine	1.371	1.249	8.9	100	-0.05
4		n-Nitrosodimethylamine	0.523	0.504	3.6	103	-0.03
5	S	2-Fluorophenol	1.278	1.110	13.1	100	-0.02
6		Aniline	2.015	1.764	12.5	98	-0.02
7	S	Phenol-d6	1.564	1.359	13.1	95	-0.01
8		2-Chlorophenol	1.521	1.340	11.9	106	-0.02
9		Benzaldehyde	0.943	0.838	11.1	97	-0.02
10	C	Phenol	1.745	1.417	18.8	95	-0.01
11		bis(2-Chloroethyl)ether	1.326	1.187	10.5	91	-0.02
12		1,3-Dichlorobenzene	1.636	1.561	4.6	112	-0.02
13	C	1,4-Dichlorobenzene	1.670	1.455	12.9	113	-0.02
14		1,2-Dichlorobenzene	1.528	1.455	4.8	110	-0.02
15		Benzyl Alcohol	1.184	1.152	2.7	119	-0.02
16		2,2'-oxybis(1-Chloropropane	1.367	1.247	8.8	109	-0.02
17		2-Methylphenol	1.193	1.010	15.3	100	-0.02
18		Hexachloroethane	0.613	0.556	9.3	109	-0.02
19	P	n-Nitroso-di-n-propylamine	0.912	0.878	3.7	116	-0.02
20		3+4-Methylphenols	1.402	1.353	3.5	122	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	110	-0.02
22		Acetophenone	0.482	0.458	5.0	113	-0.02
23	S	Nitrobenzene-d5	0.350	0.361	-3.1	122	-0.01
24		Nitrobenzene	0.362	0.324	10.5	98	-0.02
25		Isophorone	0.659	0.630	4.4	114	-0.02
26	C	2-Nitrophenol	0.188	0.200	-6.4	114	-0.02
27		2,4-Dimethylphenol	0.323	0.337	-4.3	127	-0.01
28		bis(2-Chloroethoxy)methane	0.390	0.393	-0.8	128	-0.01
29	C	2,4-Dichlorophenol	0.293	0.282	3.8	109	-0.01
30		1,2,4-Trichlorobenzene	0.315	0.292	7.3	116	-0.02
31		Naphthalene	1.130	0.944	16.5	107	-0.02
32		Benzoic acid	0.102	0.166	-62.7#	189#	-0.01
33		4-Chloroaniline	0.454	0.436	4.0	122	-0.01
34	C	Hexachlorobutadiene	0.197	0.187	5.1	112	-0.02
35		Caprolactam	0.097	0.094	3.1	123	-0.01
36	C	4-Chloro-3-methylphenol	0.367	0.339	7.6	111	-0.01
37		2-Methylnaphthalene	0.695	0.695	0.0	121	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	115	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.631	0.562	10.9	118	-0.02
40	P	Hexachlorocyclopentadiene	0.148	0.118	20.3	98	-0.03
41	S	2,4,6-Tribromophenol	0.220	0.221	-0.5	122	-0.02
42	C	2,4,6-Trichlorophenol	0.416	0.381	8.4	114	-0.02
43		2,4,5-Trichlorophenol	0.412	0.397	3.6	117	-0.02
44	S	2-Fluorobiphenyl	1.544	1.380	10.6	115	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122015\
Data File : BF083949.D
Acq On : 19 Dec 2015 12:42
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 21 02:59:21 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 18 14:23:22 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	1.885	1.558	17.3	111	-0.02
46		2-Chloronaphthalene	1.461	1.267	13.3	117	-0.02
47		2-Nitroaniline	0.396	0.395	0.3	128	-0.01
48		Acenaphthylene	2.332	2.215	5.0	120	-0.02
49		Dimethylphthalate	1.678	1.581	5.8	116	-0.02
50		2,6-Dinitrotoluene	0.366	0.320	12.6	103	-0.02
51	C	Acenaphthene	1.434	1.376	4.0	118	-0.02
52		3-Nitroaniline	0.386	0.376	2.6	122	-0.01
53	P	2,4-Dinitrophenol	0.098	0.110	-12.2	126	-0.02
54		Dibenzofuran	1.862	1.747	6.2	118	-0.02
55	P	4-Nitrophenol	0.241	0.240	0.4	108	-0.01
56		2,4-Dinitrotoluene	0.469	0.447	4.7	119	-0.01
57		Fluorene	1.603	1.475	8.0	122	-0.02
58		2,3,4,6-Tetrachlorophenol	0.305	0.278	8.9	111	-0.02
59		Diethylphthalate	1.735	1.666	4.0	119	-0.02
60		4-Chlorophenyl-phenylether	0.714	0.678	5.0	115	-0.02
61		4-Nitroaniline	0.399	0.382	4.3	127	-0.01
62		Azobenzene	1.454	1.431	1.6	116	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	114	-0.02
64		4,6-Dinitro-2-methylphenol	0.103	0.121	-17.5	115	-0.01
65	c	n-Nitrosodiphenylamine	0.728	0.645	11.4	114	-0.02
66		4-Bromophenyl-phenylether	0.245	0.233	4.9	118	-0.02
67		Hexachlorobenzene	0.257	0.248	3.5	112	-0.02
68		Atrazine	0.231	0.235	-1.7	134	-0.01
69	C	Pentachlorophenol	0.084	0.110	-31.0#	137	-0.02
70		Phenanthrene	1.208	1.124	7.0	116	-0.02
71		Anthracene	1.168	1.057	9.5	108	-0.02
72		Carbazole	1.154	1.086	5.9	122	-0.01
73		Di-n-butylphthalate	1.356	1.426	-5.2	120	-0.02
74	C	Fluoranthene	1.206	1.132	6.1	117	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	98	-0.02
76		Benzidine	0.801	0.560	30.1#	73	-0.02
77		Pyrene	1.495	1.354	9.4	114	-0.02
78	S	Terphenyl-d14	0.995	0.975	2.0	109	-0.02
79		Butylbenzylphthalate	0.666	0.675	-1.4	112	-0.02
80		Benzo(a)anthracene	1.230	1.155	6.1	106	-0.02
81		3,3'-Dichlorobenzidine	0.460	0.440	4.3	104	-0.02
82		Chrysene	1.188	1.100	7.4	106	-0.02
83		Bis(2-ethylhexyl)phthalate	0.889	0.939	-5.6	117	-0.02
84	c	Di-n-octyl phthalate	1.391	1.593	-14.5	116	-0.02
85		Indeno(1,2,3-cd)pyrene	0.888	0.930	-4.7	106	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	93	-0.03
87		Benzo(b)fluoranthene	1.458	1.531	-5.0	107	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF122015\
Data File : BF083949.D
Acq On : 19 Dec 2015 12:42
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 21 02:59:21 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 18 14:23:22 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	1.203	0.998	17.0	80	-0.02
89 C	Benzo(a)pyrene	1.204	1.127	6.4	91	-0.03
90	Dibenzo(a,h)anthracene	0.937	1.005	-7.3	105	-0.03
91	Benzo(g,h,i)perylene	0.918	1.013	-10.3	109	-0.05

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