Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100516\
Data File : BF090380.D
Acq On : 5 Oct 2016 14:41
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF100516

Quant Time: Oct 05 15:20:58 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF100516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 05 14:57:00 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	108	0.00
2		1,4-Dioxane	0.661	0.673	-1.8	110	0.00
3		Pyridine	1.842	1.935	-5.0	110	0.00
4		n-Nitrosodimethylamine	0.760	0.793	-4.3	111	0.00
5	S	2-Fluorophenol	1.340	1.346	-0.4	106	0.00
6		Aniline	2.424	2.566	-5.9	110	0.01
7	S	Phenol-d6	1.755	1.793	-2.2	112	0.00
8		2-Chlorophenol	1.475	1.469	0.4	111	0.00
9		Benzaldehyde	0.887	0.900	-1.5	111	0.00
10	C	Phenol	2.041	2.051	-0.5	114	0.00
11		bis(2-Chloroethyl)ether	1.562	1.646	-5.4	107	0.00
12		1,3-Dichlorobenzene	1.470	1.465	0.3	110	0.00
13	C	1,4-Dichlorobenzene	1.493	1.629	-9.1	111	0.00
14		1,2-Dichlorobenzene	1.439	1.400	2.7	113	0.01
15		Benzyl Alcohol	1.334	1.370	-2.7	116	0.00
16		2,2'-oxybis(1-Chloropropane	2.650	2.784	-5.1	106	0.01
17		2-Methylphenol	1.213	1.217	-0.3	109	0.00
18		Hexachloroethane	0.589	0.617	-4.8	111	0.00
19	P	n-Nitroso-di-n-propylamine	1.287	1.390	-8.0	111	0.00
20		3+4-Methylphenols	1.570	1.704	-8.5	111	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
22		Acetophenone	0.474	0.466	1.7	109	0.00
23	S	Nitrobenzene-d5	0.411	0.429	-4.4	107	0.00
24		Nitrobenzene	0.408	0.407	0.2	111	0.00
25		Isophorone	0.752	0.788	-4.8	114	0.00
26	C	2-Nitrophenol	0.176	0.170	3.4	106	0.00
27		2,4-Dimethylphenol	0.342	0.320	6.4	108	0.00
28		bis(2-Chloroethoxy)methane	0.445	0.501	-12.6	110	0.00
29	C	2,4-Dichlorophenol	0.260	0.255	1.9	113	0.00
30		1,2,4-Trichlorobenzene	0.268	0.286	-6.7	109	0.00
31		Naphthalene	0.965	0.914	5.3	106	0.00
32		Benzoic acid	0.238	0.267	-12.2	123	0.01
33		4-Chloroaniline	0.399	0.430	-7.8	107	0.00
34	C	Hexachlorobutadiene	0.150	0.160	-6.7	113	0.00
35		Caprolactam	0.087	0.090	-3.4	109	0.00
36	C	4-Chloro-3-methylphenol	0.325	0.340	-4.6	106	0.00
37		2-Methylnaphthalene	0.584	0.645	-10.4	109	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
39		1,2,4,5-Tetrachlorobenzene	0.527	0.525	0.4	115	0.00
40	P	Hexachlorocyclopentadiene	0.289	0.317	-9.7	114	0.00
41	S	2,4,6-Tribromophenol	0.163	0.179	-9.8	111	0.00
42	C	2,4,6-Trichlorophenol	0.364	0.416	-14.3	108	0.00
43		2,4,5-Trichlorophenol	0.344	0.332	3.5	108	0.00
44	S	2-Fluorobiphenyl	1.200	1.336	-11.3	108	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100516\
Data File : BF090380.D
Acq On : 5 Oct 2016 14:41
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF100516

Quant Time: Oct 05 15:20:58 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF100516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 05 14:57:00 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	1.482	1.585	-7.0	112	0.00
46		2-Chloronaphthalene	1.095	1.258	-14.9	110	0.00
47		2-Nitroaniline	0.432	0.484	-12.0	114	0.00
48		Acenaphthylene	1.807	1.750	3.2	108	0.00
49		Dimethylphthalate	1.281	1.405	-9.7	111	0.00
50		2,6-Dinitrotoluene	0.284	0.315	-10.9	108	0.00
51	C	Acenaphthene	1.127	1.141	-1.2	112	0.00
52		3-Nitroaniline	0.332	0.398	-19.9	117	0.00
53	P	2,4-Dinitrophenol	0.127	0.143	-12.6	124	0.01
54		Dibenzofuran	1.458	1.437	1.4	112	0.00
55	P	4-Nitrophenol	0.291	0.332	-14.1	108	0.00
56		2,4-Dinitrotoluene	0.378	0.372	1.6	103	0.00
57		Fluorene	1.225	1.296	-5.8	114	0.00
58		2,3,4,6-Tetrachlorophenol	0.257	0.248	3.5	110	0.00
59		Diethylphthalate	1.325	1.345	-1.5	110	0.00
60		4-Chlorophenyl-phenylether	0.557	0.646	-16.0	113	0.00
61		4-Nitroaniline	0.335	0.373	-11.3	122	0.00
62		Azobenzene	1.531	1.714	-12.0	108	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
64		4,6-Dinitro-2-methylphenol	0.121	0.128	-5.8	108	0.00
65	c	n-Nitrosodiphenylamine	0.670	0.672	-0.3	108	0.00
66		4-Bromophenyl-phenylether	0.196	0.214	-9.2	109	0.00
67		Hexachlorobenzene	0.218	0.242	-11.0	112	0.00
68		Atrazine	0.165	0.166	-0.6	108	0.00
69	C	Pentachlorophenol	0.130	0.127	2.3	112	0.00
70		Phenanthrene	1.024	1.078	-5.3	112	0.00
71		Anthracene	1.046	1.017	2.8	111	0.00
72		Carbazole	0.972	1.075	-10.6	110	0.00
73		Di-n-butylphthalate	1.182	1.319	-11.6	111	0.00
74	C	Fluoranthene	1.005	0.928	7.7	111	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	110	0.00
76		Benzidine	0.526	0.503	4.4	110	0.00
77		Pyrene	1.334	1.258	5.7	111	0.00
78	S	Terphenyl-d14	0.890	0.868	2.5	111	0.00
79		Butylbenzylphthalate	0.684	0.665	2.8	111	0.00
80		Benzo(a)anthracene	1.122	1.158	-3.2	115	0.00
81		3,3'-Dichlorobenzidine	0.407	0.375	7.9	115	0.00
82		Chrysene	1.033	1.033	0.0	112	0.00
83		Bis(2-ethylhexyl)phthalate	0.973	1.035	-6.4	115	0.00
84	c	Di-n-octyl phthalate	1.442	1.486	-3.1	115	0.00
85		Indeno(1,2,3-cd)pyrene	0.736	0.769	-4.5	123	0.01
86	I	Perylene-d12	1.000	1.000	0.0	117	0.00
87		Benzo(b)fluoranthene	1.274	1.382	-8.5	116	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF100516\
Data File : BF090380.D
Acq On : 5 Oct 2016 14:41
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF100516

Quant Time: Oct 05 15:20:58 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF100516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 05 14:57:00 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	1.197	1.111	7.2	116	0.00
89 C	Benzo(a)pyrene	1.098	1.133	-3.2	117	0.00
90	Dibenzo(a,h)anthracene	0.934	0.980	-4.9	122	0.00
91	Benzo(g,h,i)perylene	0.896	0.937	-4.6	122	0.00

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