Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF083116\
Data File : BF089936.D
Acq On : 31 Aug 2016 5:57
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF083116

Quant Time: Aug 31 11:57:20 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 31 11:55:18 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	103	0.00
2		1,4-Dioxane	0.592	0.564	4.7	105	0.00
3		Pyridine	1.579	1.611	-2.0	101	0.00
4		n-Nitrosodimethylamine	0.779	0.818	-5.0	105	0.00
5	S	2-Fluorophenol	1.250	1.300	-4.0	103	0.00
6		Aniline	2.072	1.947	6.0	102	0.00
7	S	Phenol-d6	1.518	1.563	-3.0	105	0.01
8		2-Chlorophenol	1.356	1.397	-3.0	105	0.00
9		Benzaldehyde	0.980	0.978	0.2	103	0.00
10	C	Phenol	1.766	1.813	-2.7	109	0.00
11		bis(2-Chloroethyl)ether	1.336	1.273	4.7	104	0.00
12		1,3-Dichlorobenzene	1.514	1.493	1.4	105	0.00
13	C	1,4-Dichlorobenzene	1.549	1.483	4.3	103	0.00
14		1,2-Dichlorobenzene	1.389	1.453	-4.6	105	0.00
15		Benzyl Alcohol	0.951	0.935	1.7	101	0.00
16		2,2'-oxybis(1-Chloropropane	2.547	2.713	-6.5	108	0.00
17		2-Methylphenol	1.083	1.110	-2.5	106	0.00
18		Hexachloroethane	0.531	0.534	-0.6	105	0.00
19	P	n-Nitroso-di-n-propylamine	0.989	1.041	-5.3	104	0.00
20		3+4-Methylphenols	1.312	1.159	11.7	103	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
22		Acetophenone	0.449	0.445	0.9	108	0.00
23	S	Nitrobenzene-d5	0.345	0.334	3.2	103	0.00
24		Nitrobenzene	0.367	0.340	7.4	105	0.00
25		Isophorone	0.701	0.685	2.3	106	0.00
26	C	2-Nitrophenol	0.175	0.173	1.1	105	0.00
27		2,4-Dimethylphenol	0.324	0.312	3.7	112	0.00
28		bis(2-Chloroethoxy)methane	0.423	0.443	-4.7	107	0.00
29	C	2,4-Dichlorophenol	0.291	0.271	6.9	105	0.00
30		1,2,4-Trichlorobenzene	0.318	0.321	-0.9	104	0.00
31		Naphthalene	0.996	0.887	10.9	102	0.00
32		Benzoic acid	0.218	0.240	-10.1	117	0.01
33		4-Chloroaniline	0.402	0.409	-1.7	106	0.00
34	C	Hexachlorobutadiene	0.186	0.178	4.3	104	0.00
35		Caprolactam	0.092	0.088	4.3	97	0.00
36	C	4-Chloro-3-methylphenol	0.308	0.322	-4.5	111	0.00
37		2-Methylnaphthalene	0.658	0.669	-1.7	106	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	105	0.00
39		1,2,4,5-Tetrachlorobenzene	0.586	0.524	10.6	106	0.00
40	P	Hexachlorocyclopentadiene	0.300	0.315	-5.0	107	-0.01
41	S	2,4,6-Tribromophenol	0.197	0.212	-7.6	112	0.00
42	C	2,4,6-Trichlorophenol	0.406	0.438	-7.9	107	0.00
43		2,4,5-Trichlorophenol	0.390	0.372	4.6	108	0.00
44	S	2-Fluorobiphenyl	1.218	1.167	4.2	104	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF083116\
Data File : BF089936.D
Acq On : 31 Aug 2016 5:57
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF083116

Quant Time: Aug 31 11:57:20 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 31 11:55:18 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	1.610	1.463	9.1	105	0.00
46		2-Chloronaphthalene	1.139	1.097	3.7	105	0.00
47		2-Nitroaniline	0.372	0.382	-2.7	111	0.00
48		Acenaphthylene	1.884	1.784	5.3	102	0.00
49		Dimethylphthalate	1.446	1.399	3.3	115	0.00
50		2,6-Dinitrotoluene	0.322	0.308	4.3	103	0.00
51	C	Acenaphthene	1.193	1.120	6.1	103	0.00
52		3-Nitroaniline	0.367	0.390	-6.3	124	0.01
53	P	2,4-Dinitrophenol	0.130	0.160	-23.1	116	0.00
54		Dibenzofuran	1.599	1.562	2.3	106	0.00
55	P	4-Nitrophenol	0.277	0.328	-18.4	115	0.00
56		2,4-Dinitrotoluene	0.408	0.442	-8.3	102	0.00
57		Fluorene	1.169	1.240	-6.1	107	0.00
58		2,3,4,6-Tetrachlorophenol	0.334	0.321	3.9	105	0.00
59		Diethylphthalate	1.360	1.331	2.1	109	0.00
60		4-Chlorophenyl-phenylether	0.550	0.571	-3.8	108	0.00
61		4-Nitroaniline	0.360	0.394	-9.4	106	0.00
62		Azobenzene	1.358	1.317	3.0	110	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	109	0.00
64		4,6-Dinitro-2-methylphenol	0.128	0.116	9.4	108	0.00
65	c	n-Nitrosodiphenylamine	0.601	0.539	10.3	104	0.00
66		4-Bromophenyl-phenylether	0.201	0.190	5.5	107	0.00
67		Hexachlorobenzene	0.230	0.210	8.7	112	0.00
68		Atrazine	0.201	0.205	-2.0	120	0.01
69	C	Pentachlorophenol	0.124	0.137	-10.5	108	0.00
70		Phenanthrene	0.996	0.951	4.5	106	0.00
71		Anthracene	1.010	0.913	9.6	112	0.00
72		Carbazole	0.947	0.870	8.1	105	0.00
73		Di-n-butylphthalate	1.104	1.100	0.4	116	0.00
74	C	Fluoranthene	1.044	1.071	-2.6	110	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
76		Benzidine	0.758	0.778	-2.6	106	0.00
77		Pyrene	1.410	1.336	5.2	108	0.00
78	S	Terphenyl-d14	0.810	0.732	9.6	111	0.00
79		Butylbenzylphthalate	0.569	0.580	-1.9	108	0.00
80		Benzo(a)anthracene	1.225	1.202	1.9	104	0.00
81		3,3'-Dichlorobenzidine	0.442	0.441	0.2	106	0.00
82		Chrysene	1.144	1.120	2.1	107	0.00
83		Bis(2-ethylhexyl)phthalate	0.787	0.800	-1.7	105	-0.01
84	c	Di-n-octyl phthalate	1.270	1.454	-14.5	110	0.00
85		Indeno(1,2,3-cd)pyrene	0.846	0.920	-8.7	113	0.01
86	I	Perylene-d12	1.000	1.000	0.0	110	0.00
87		Benzo(b)fluoranthene	1.253	1.389	-10.9	108	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF083116\
Data File : BF089936.D
Acq On : 31 Aug 2016 5:57
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF083116

Quant Time: Aug 31 11:57:20 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF083116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 31 11:55:18 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	1.063	0.867	18.4	107	0.00
89 C	Benzo(a)pyrene	1.073	1.086	-1.2	110	0.00
90	Dibenzo(a,h)anthracene	0.787	0.798	-1.4	112	0.00
91	Benzo(g,h,i)perylene	0.792	0.790	0.3	112	0.00

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