Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110715\
Data File : BF082752.D
Acq On : 6 Nov 2015 10:20
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040

Quant Time: Nov 06 17:15:27 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 06 14:39:58 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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	121	0.00
2		1,4-Dioxane	40.000	37.720	5.7	112	0.00
3		Pyridine	40.000	38.010	5.0	111	0.00
4		n-Nitrosodimethylamine	40.000	39.213	2.0	121	0.02
5	S	2-Fluorophenol	80.000	69.675	12.9	104	0.00
6		Aniline	40.000	39.218	2.0	125	0.01
7	S	Phenol-d6	80.000	76.173	4.8	119	0.01
8		2-Chlorophenol	40.000	36.533	8.7	111	0.00
9		Benzaldehyde	40.000	39.065	2.3	112	0.00
10	C	Phenol	40.000	35.797	10.5	103	0.01
11		bis(2-Chloroethyl)ether	40.000	34.769	13.1	100	0.00
12		1,3-Dichlorobenzene	40.000	33.267	16.8	99	0.23
13	C	1,4-Dichlorobenzene	40.000	39.381	1.5	131	0.01
14		1,2-Dichlorobenzene	40.000	35.176	12.1	106	0.00
15		Benzyl Alcohol	40.000	40.487	-1.2	127	0.01
16		2,2'-oxybis(1-Chloropropane	40.000	38.174	4.6	119	0.00
17		2-Methylphenol	40.000	38.488	3.8	115	0.01
18		Hexachloroethane	40.000	38.718	3.2	125	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.523	1.2	122	0.02
20		3+4-Methylphenols	40.000	40.755	-1.9	137	0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	129	0.00
22		Acetophenone	40.000	37.278	6.8	132	0.01
23	S	Nitrobenzene-d5	80.000	71.848	10.2	117	0.01
24		Nitrobenzene	40.000	35.130	12.2	118	0.00
25		Isophorone	40.000	36.822	7.9	124	0.01
26	C	2-Nitrophenol	40.000	35.044	12.4	104	0.00
27		2,4-Dimethylphenol	40.000	36.682	8.3	124	0.01
28		bis(2-Chloroethoxy)methane	40.000	33.693	15.8	106	0.00
29	C	2,4-Dichlorophenol	40.000	35.875	10.3	120	0.00
30		1,2,4-Trichlorobenzene	40.000	37.424	6.4	122	0.00
31		Naphthalene	40.000	36.231	9.4	122	0.01
32		Benzoic acid	40.000	39.447	1.4	125	0.03
33		4-Chloroaniline	40.000	36.181	9.5	118	0.01
34	C	Hexachlorobutadiene	40.000	39.021	2.4	129	0.00
35		Caprolactam	40.000	37.346	6.6	116	0.03
36	C	4-Chloro-3-methylphenol	40.000	35.178	12.1	117	0.00
37		2-Methylnaphthalene	40.000	34.309	14.2	116	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	124	0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	37.797	5.5	108	0.00
40	P	Hexachlorocyclopentadiene	40.000	40.674	-1.7	120	0.00
41	S	2,4,6-Tribromophenol	80.000	77.061	3.7	125	0.00
42	C	2,4,6-Trichlorophenol	40.000	39.885	0.3	128	0.01
43		2,4,5-Trichlorophenol	40.000	38.766	3.1	115	0.01
44	S	2-Fluorobiphenyl	80.000	78.815	1.5	131	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110715\
Data File : BF082752.D
Acq On : 6 Nov 2015 10:20
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040

Quant Time: Nov 06 17:15:27 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 06 14:39:58 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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	35.454	11.4	102	0.00
46		2-Chloronaphthalene	40.000	36.002	10.0	105	0.00
47		2-Nitroaniline	40.000	38.138	4.7	107	0.01
48		Acenaphthylene	40.000	37.841	5.4	121	0.00
49		Dimethylphthalate	40.000	40.212	-0.5	137	0.01
50		2,6-Dinitrotoluene	40.000	42.163	-5.4	146	0.01
51	C	Acenaphthene	40.000	39.921	0.2	128	0.01
52		3-Nitroaniline	40.000	34.312	14.2	95	0.01
53	P	2,4-Dinitrophenol	40.000	37.453	6.4	100	0.01
54		Dibenzofuran	40.000	37.767	5.6	122	0.00
55	P	4-Nitrophenol	40.000	35.778	10.6	112	0.01
56		2,4-Dinitrotoluene	40.000	41.722	-4.3	143	0.01
57		Fluorene	40.000	36.166	9.6	114	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	42.055	-5.1	140	-0.03
59		Diethylphthalate	40.000	41.027	-2.6	129	0.01
60		4-Chlorophenyl-phenylether	40.000	40.896	-2.2	133	0.01
61		4-Nitroaniline	40.000	36.779	8.1	118	0.01
62		Azobenzene	40.000	42.138	-5.3	131	0.01
63	I	Phenanthrene-d10	20.000	20.000	0.0	125	0.00
64		4,6-Dinitro-2-methylphenol	40.000	43.536	-8.8	139	0.01
65	c	n-Nitrosodiphenylamine	40.000	33.323	16.7	104	0.00
66		4-Bromophenyl-phenylether	40.000	39.989	0.0	137	0.01
67		Hexachlorobenzene	40.000	40.520	-1.3	139	0.01
68		Atrazine	40.000	39.507	1.2	131	0.01
69	C	Pentachlorophenol	40.000	39.121	2.2	111	0.00
70		Phenanthrene	40.000	35.585	11.0	111	0.00
71		Anthracene	40.000	37.763	5.6	127	0.01
72		Carbazole	40.000	33.162	17.1	102	0.00
73		Di-n-butylphthalate	40.000	38.629	3.4	123	0.00
74	C	Fluoranthene	40.000	34.264	14.3	106	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	117	0.01
76		Benzidine	40.000	39.489	1.3	110	0.00
77		Pyrene	40.000	36.265	9.3	110	0.01
78	S	Terphenyl-d14	80.000	70.651	11.7	102	0.00
79		Butylbenzylphthalate	40.000	45.182	-13.0	141	0.01
80		Benzo(a)anthracene	40.000	37.838	5.4	117	0.00
81		3,3'-Dichlorobenzidine	40.000	36.283	9.3	107	0.01
82		Chrysene	40.000	39.564	1.1	126	0.01
83		Bis(2-ethylhexyl)phthalate	40.000	44.568	-11.4	138	0.01
84	c	Di-n-octyl phthalate	40.000	40.329	-0.8	120	0.01
85		Indeno(1,2,3-cd)pyrene	40.000	37.673	5.8	117	0.03
86	I	Perylene-d12	20.000	20.000	0.0	122	0.02
87		Benzo(b)fluoranthene	40.000	36.528	8.7	123	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF110715\
Data File : BF082752.D
Acq On : 6 Nov 2015 10:20
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040

Quant Time: Nov 06 17:15:27 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF110715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Nov 06 14:39:58 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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	41.177	-2.9	120	-0.01
89 C	Benzo(a)pyrene	40.000	37.553	6.1	117	0.02
90	Dibenzo(a,h)anthracene	40.000	37.567	6.1	117	0.03
91	Benzo(g,h,i)perylene	40.000	36.489	8.8	116	0.03

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