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	Amount	Calc.	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	20.000	20.000	0.0	111	-0.02
2		1,4-Dioxane	40.000	33.890	15.3	95	-0.05
3		Pyridine	40.000	36.458	8.9	100	-0.05
4		n-Nitrosodimethylamine	40.000	38.524	3.7	103	-0.03
5	S	2-Fluorophenol	80.000	69.464	13.2	100	-0.02
6		Aniline	40.000	35.017	12.5	98	-0.02
7	S	Phenol-d6	80.000	69.481	13.1	95	-0.01
8		2-Chlorophenol	40.000	35.244	11.9	106	-0.02
9		Benzaldehyde	40.000	35.576	11.1	97	-0.02
10	C	Phenol	40.000	32.468	18.8	95	-0.01
11		bis(2-Chloroethyl)ether	40.000	35.812	10.5	91	-0.02
12		1,3-Dichlorobenzene	40.000	38.163	4.6	112	-0.02
13	C	1,4-Dichlorobenzene	40.000	34.844	12.9	113	-0.02
14		1,2-Dichlorobenzene	40.000	43.094	-7.7	110	-0.02
15		Benzyl Alcohol	40.000	38.931	2.7	119	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	36.466	8.8	109	-0.02
17		2-Methylphenol	40.000	33.846	15.4	100	-0.02
18		Hexachloroethane	40.000	36.296	9.3	109	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	38.510	3.7	116	-0.02
20		3+4-Methylphenols	40.000	38.615	3.5	122	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	110	-0.02
22		Acetophenone	40.000	38.046	4.9	113	-0.02
23	S	Nitrobenzene-d5	80.000	82.303	-2.9	122	-0.01
24		Nitrobenzene	40.000	35.794	10.5	98	-0.02
25		Isophorone	40.000	38.199	4.5	114	-0.02
26	C	2-Nitrophenol	40.000	42.573	-6.4	114	-0.02
27		2,4-Dimethylphenol	40.000	41.682	-4.2	127	-0.01
28		bis(2-Chloroethoxy)methane	40.000	40.305	-0.8	128	-0.01
29	C	2,4-Dichlorophenol	40.000	38.532	3.7	109	-0.01
30		1,2,4-Trichlorobenzene	40.000	37.104	7.2	116	-0.02
31		Naphthalene	40.000	37.080	7.3	107	-0.02
32		Benzoic acid	40.000	53.525	-33.8#	189	-0.01
33		4-Chloroaniline	40.000	38.420	3.9	122	-0.01
34	C	Hexachlorobutadiene	40.000	38.055	4.9	112	-0.02
35		Caprolactam	40.000	38.663	3.3	123	-0.01
36	C	4-Chloro-3-methylphenol	40.000	40.803	-2.0	111	-0.01
37		2-Methylnaphthalene	40.000	39.981	0.0	121	-0.02
38	I	Acenaphthene-d10	20.000	20.000	0.0	115	-0.02
39		1,2,4,5-Tetrachlorobenzene	40.000	35.672	10.8	118	-0.02
40	P	Hexachlorocyclopentadiene	40.000	34.650	13.4	98	-0.03
41	S	2,4,6-Tribromophenol	80.000	80.421	-0.5	122	-0.02
42	C	2,4,6-Trichlorophenol	40.000	36.580	8.6	114	-0.02
43		2,4,5-Trichlorophenol	40.000	38.523	3.7	117	-0.02
44	S	2-Fluorobiphenyl	80.000	76.242	4.7	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	Amount	Calc.	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	40.000	36.834	7.9	111	-0.02
46		2-Chloronaphthalene	40.000	38.289	4.3	117	-0.02
47		2-Nitroaniline	40.000	39.945	0.1	128	-0.01
48		Acenaphthylene	40.000	38.000	5.0	120	-0.02
49		Dimethylphthalate	40.000	37.672	5.8	116	-0.02
50		2,6-Dinitrotoluene	40.000	34.998	12.5	103	-0.02
51	C	Acenaphthene	40.000	38.404	4.0	118	-0.02
52		3-Nitroaniline	40.000	39.003	2.5	122	-0.01
53	P	2,4-Dinitrophenol	40.000	39.466	1.3	126	-0.02
54		Dibenzofuran	40.000	37.532	6.2	118	-0.02
55	P	4-Nitrophenol	40.000	39.826	0.4	108	-0.01
56		2,4-Dinitrotoluene	40.000	38.135	4.7	119	-0.01
57		Fluorene	40.000	41.434	-3.6	122	-0.02
58		2,3,4,6-Tetrachlorophenol	40.000	36.451	8.9	111	-0.02
59		Diethylphthalate	40.000	38.406	4.0	119	-0.02
60		4-Chlorophenyl-phenylether	40.000	38.009	5.0	115	-0.02
61		4-Nitroaniline	40.000	38.313	4.2	127	-0.01
62		Azobenzene	40.000	39.366	1.6	116	-0.02
63	I	Phenanthrene-d10	20.000	20.000	0.0	114	-0.02
64		4,6-Dinitro-2-methylphenol	40.000	40.052	-0.1	115	-0.01
65	c	n-Nitrosodiphenylamine	40.000	35.470	11.3	114	-0.02
66		4-Bromophenyl-phenylether	40.000	38.103	4.7	118	-0.02
67		Hexachlorobenzene	40.000	38.525	3.7	112	-0.02
68		Atrazine	40.000	40.790	-2.0	134	-0.01
69	C	Pentachlorophenol	40.000	42.575	-6.4	137	-0.02
70		Phenanthrene	40.000	37.195	7.0	116	-0.02
71		Anthracene	40.000	36.206	9.5	108	-0.02
72		Carbazole	40.000	37.626	5.9	122	-0.01
73		Di-n-butylphthalate	40.000	42.065	-5.2	120	-0.02
74	C	Fluoranthene	40.000	37.559	6.1	117	-0.02
75	I	Chrysene-d12	20.000	20.000	0.0	98	-0.02
76		Benzidine	40.000	27.972	30.1#	73	-0.02
77		Pyrene	40.000	42.759	-6.9	114	-0.02
78	S	Terphenyl-d14	80.000	91.699	-14.6	109	-0.02
79		Butylbenzylphthalate	40.000	40.533	-1.3	112	-0.02
80		Benzo(a)anthracene	40.000	37.566	6.1	106	-0.02
81		3,3'-Dichlorobenzidine	40.000	38.326	4.2	104	-0.02
82		Chrysene	40.000	37.041	7.4	106	-0.02
83		Bis(2-ethylhexyl)phthalate	40.000	42.240	-5.6	117	-0.02
84	c	Di-n-octyl phthalate	40.000	45.793	-14.5	116	-0.02
85		Indeno(1,2,3-cd)pyrene	40.000	41.861	-4.7	106	-0.03
86	I	Perylene-d12	20.000	20.000	0.0	93	-0.03
87		Benzo(b)fluoranthene	40.000	42.001	-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	Amount	Calc.	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	40.000	33.198	17.0	80	-0.02
89 C	Benzo(a)pyrene	40.000	37.430	6.4	91	-0.03
90	Dibenzo(a,h)anthracene	40.000	42.923	-7.3	105	-0.03
91	Benzo(g,h,i)perylene	40.000	44.141	-10.4	109	-0.05

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