Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121015\
Data File : BF083684.D
Acq On : 10 Dec 2015 19:01
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 17 11:22:12 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 17 09:56:43 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	81	-0.09
2		1,4-Dioxane	40.000	42.208	-5.5	89	-0.09
3		Pyridine	40.000	40.022	-0.1	82	-0.09
4		n-Nitrosodimethylamine	40.000	34.968	12.6	66	-0.08
5	S	2-Fluorophenol	80.000	82.888	-3.6	80	-0.08
6		Aniline	40.000	39.695	0.8	75	-0.08
7	S	Phenol-d6	80.000	80.841	-1.1	79	-0.09
8		2-Chlorophenol	40.000	40.696	-1.7	81	-0.09
9		Benzaldehyde	40.000	42.073	-5.2	83	-0.08
10	C	Phenol	40.000	40.723	-1.8	79	-0.08
11		bis(2-Chloroethyl)ether	40.000	40.868	-2.2	81	-0.08
12		1,3-Dichlorobenzene	40.000	40.729	-1.8	81	-0.08
13	C	1,4-Dichlorobenzene	40.000	39.979	0.1	81	-0.09
14		1,2-Dichlorobenzene	40.000	40.487	-1.2	80	-0.09
15		Benzyl Alcohol	40.000	39.479	1.3	74	-0.08
16		2,2'-oxybis(1-Chloropropane	40.000	39.047	2.4	77	-0.08
17		2-Methylphenol	40.000	39.926	0.2	77	-0.08
18		Hexachloroethane	40.000	39.779	0.6	78	-0.10
19	P	n-Nitroso-di-n-propylamine	40.000	39.413	1.5	79	-0.09
20		3+4-Methylphenols	40.000	37.688	5.8	76	-0.08
21	I	Naphthalene-d8	20.000	20.000	0.0	76	-0.10
22		Acetophenone	40.000	40.733	-1.8	76	-0.08
23	S	Nitrobenzene-d5	80.000	83.087	-3.9	79	-0.09
24		Nitrobenzene	40.000	41.357	-3.4	80	-0.09
25		Isophorone	40.000	41.859	-4.6	79	-0.09
26	C	2-Nitrophenol	40.000	43.183	-8.0	78	-0.09
27		2,4-Dimethylphenol	40.000	43.123	-7.8	81	-0.09
28		bis(2-Chloroethoxy)methane	40.000	41.537	-3.8	78	-0.09
29	C	2,4-Dichlorophenol	40.000	41.767	-4.4	77	-0.08
30		1,2,4-Trichlorobenzene	40.000	41.166	-2.9	78	-0.09
31		Naphthalene	40.000	41.825	-4.6	81	-0.09
32		Benzoic acid	40.000	31.748	20.6	62	-0.07
33		4-Chloroaniline	40.000	35.950	10.1	66	-0.07
34	C	Hexachlorobutadiene	40.000	42.007	-5.0	80	-0.10
35		Caprolactam	40.000	44.347	-10.9	80	-0.09
36	C	4-Chloro-3-methylphenol	40.000	43.413	-8.5	81	-0.08
37		2-Methylnaphthalene	40.000	42.667	-6.7	82	-0.09
38	I	Acenaphthene-d10	20.000	20.000	0.0	78	-0.09
39		1,2,4,5-Tetrachlorobenzene	40.000	41.290	-3.2	80	-0.09
40	P	Hexachlorocyclopentadiene	40.000	25.190	37.0#	31	-0.10
41	S	2,4,6-Tribromophenol	80.000	85.014	-6.3	81	-0.10
42	C	2,4,6-Trichlorophenol	40.000	39.692	0.8	75	-0.08
43		2,4,5-Trichlorophenol	40.000	38.156	4.6	71	-0.07
44	S	2-Fluorobiphenyl	80.000	81.259	-1.6	81	-0.09

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121015\
Data File : BF083684.D
Acq On : 10 Dec 2015 19:01
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 17 11:22:12 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 17 09:56:43 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	41.811	-4.5	82	-0.09
46		2-Chloronaphthalene	40.000	41.730	-4.3	81	-0.09
47		2-Nitroaniline	40.000	42.128	-5.3	80	-0.08
48		Acenaphthylene	40.000	41.298	-3.2	81	-0.09
49		Dimethylphthalate	40.000	40.460	-1.2	80	-0.09
50		2,6-Dinitrotoluene	40.000	41.868	-4.7	80	-0.09
51	C	Acenaphthene	40.000	41.412	-3.5	81	-0.09
52		3-Nitroaniline	40.000	42.250	-5.6	78	-0.09
53	P	2,4-Dinitrophenol	40.000	41.097	-2.7	82	-0.05
54		Dibenzofuran	40.000	41.948	-4.9	81	-0.09
55	P	4-Nitrophenol	40.000	31.669	20.8	61	-0.03
56		2,4-Dinitrotoluene	40.000	43.673	-9.2	80	-0.08
57		Fluorene	40.000	41.358	-3.4	82	-0.09
58		2,3,4,6-Tetrachlorophenol	40.000	39.280	1.8	76	-0.08
59		Diethylphthalate	40.000	40.408	-1.0	80	-0.09
60		4-Chlorophenyl-phenylether	40.000	41.660	-4.1	80	-0.10
61		4-Nitroaniline	40.000	44.807	-12.0	82	-0.08
62		Azobenzene	40.000	43.879	-9.7	84	-0.09
63	I	Phenanthrene-d10	20.000	20.000	0.0	79	-0.09
64		4,6-Dinitro-2-methylphenol	40.000	39.388	1.5	77	-0.07
65	c	n-Nitrosodiphenylamine	40.000	41.174	-2.9	81	-0.09
66		4-Bromophenyl-phenylether	40.000	42.009	-5.0	82	-0.09
67		Hexachlorobenzene	40.000	41.299	-3.2	81	-0.09
68		Atrazine	40.000	41.793	-4.5	79	-0.10
69	C	Pentachlorophenol	40.000	24.716	38.2#	44	-0.06
70		Phenanthrene	40.000	41.970	-4.9	83	-0.09
71		Anthracene	40.000	42.741	-6.9	84	-0.09
72		Carbazole	40.000	43.832	-9.6	87	-0.09
73		Di-n-butylphthalate	40.000	42.254	-5.6	82	-0.10
74	C	Fluoranthene	40.000	43.410	-8.5	85	-0.09
75	I	Chrysene-d12	20.000	20.000	0.0	85	-0.08
76		Benzidine	40.000	39.305	1.7	78	-0.08
77		Pyrene	40.000	40.711	-1.8	86	-0.09
78	S	Terphenyl-d14	80.000	80.922	-1.2	88	-0.10
79		Butylbenzylphthalate	40.000	41.729	-4.3	86	-0.09
80		Benzo(a)anthracene	40.000	40.552	-1.4	87	-0.08
81		3,3'-Dichlorobenzidine	40.000	44.676	-11.7	91	-0.08
82		Chrysene	40.000	42.826	-7.1	91	-0.08
83		Bis(2-ethylhexyl)phthalate	40.000	41.224	-3.1	84	-0.08
84	c	Di-n-octyl phthalate	40.000	42.715	-6.8	89	-0.08
85		Indeno(1,2,3-cd)pyrene	40.000	44.335	-10.8	100	-0.09
86	I	Perylene-d12	20.000	20.000	0.0	92	-0.06
87		Benzo(b)fluoranthene	40.000	44.561	-11.4	104	-0.07

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF121015\
Data File : BF083684.D
Acq On : 10 Dec 2015 19:01
Operator : UM/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 17 11:22:12 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF120815.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Dec 17 09:56:43 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	39.624	0.9	89	-0.07
89 C	Benzo(a)pyrene	40.000	42.091	-5.2	98	-0.07
90	Dibenzo(a,h)anthracene	40.000	42.951	-7.4	103	-0.10
91	Benzo(g,h,i)perylene	40.000	39.422	1.4	96	-0.09

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