Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF041415\
Data File : BF078527.D
Acq On : 15 Apr 2015 6:18
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 15 05:57:13 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 15 01:10:59 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	103	-0.08
2		1,4-Dioxane	40.000	51.396	-28.5#	133	-0.15
3		Pyridine	40.000	44.718	-11.8	110	-0.21
4		n-Nitrosodimethylamine	40.000	42.967	-7.4	114	-0.20
5	S	2-Fluorophenol	80.000	79.080	1.2	103	-0.12
6		Aniline	40.000	44.857	-12.1	119	-0.08
7	S	Phenol-d6	80.000	83.482	-4.4	110	-0.07
8		2-Chlorophenol	40.000	41.513	-3.8	108	-0.09
9		Benzaldehyde	40.000	37.053	7.4	94	-0.10
10	C	Phenol	40.000	40.860	-2.1	107	-0.07
11		bis(2-Chloroethyl)ether	40.000	41.429	-3.6	105	-0.08
12		1,3-Dichlorobenzene	40.000	39.339	1.7	105	-0.09
13	C	1,4-Dichlorobenzene	40.000	38.830	2.9	104	-0.09
14		1,2-Dichlorobenzene	40.000	39.369	1.6	105	-0.09
15		Benzyl Alcohol	40.000	48.697	-21.7	128	-0.07
16		2,2'-oxybis(1-Chloropropane	40.000	39.854	0.4	105	-0.08
17		2-Methylphenol	40.000	42.992	-7.5	110	-0.07
18		Hexachloroethane	40.000	35.849	10.4	94	-0.09
19	P	n-Nitroso-di-n-propylamine	40.000	42.820	-7.1	111	-0.08
20		3+4-Methylphenols	40.000	45.347	-13.4	116	-0.07
21	I	Naphthalene-d8	20.000	20.000	0.0	119	-0.08
22		Acetophenone	40.000	40.120	-0.3	118	-0.08
23	S	Nitrobenzene-d5	80.000	78.754	1.6	116	-0.07
24		Nitrobenzene	40.000	38.330	4.2	115	-0.09
25		Isophorone	40.000	41.290	-3.2	117	-0.08
26	C	2-Nitrophenol	40.000	34.383	14.0	93	-0.08
27		2,4-Dimethylphenol	40.000	38.531	3.7	111	-0.07
28		bis(2-Chloroethoxy)methane	40.000	37.331	6.7	108	-0.07
29	C	2,4-Dichlorophenol	40.000	39.470	1.3	115	-0.08
30		1,2,4-Trichlorobenzene	40.000	35.259	11.9	106	-0.09
31		Naphthalene	40.000	37.245	6.9	111	-0.09
32		Benzoic acid	40.000	47.215	-18.0	144	-0.05
33		4-Chloroaniline	40.000	40.732	-1.8	115	-0.08
34	C	Hexachlorobutadiene	40.000	36.293	9.3	107	-0.09
35		Caprolactam	40.000	49.273	-23.2	136	-0.04
36	C	4-Chloro-3-methylphenol	40.000	42.205	-5.5	120	-0.07
37		2-Methylnaphthalene	40.000	38.254	4.4	113	-0.09
38	I	Acenaphthene-d10	20.000	20.000	0.0	122	-0.09
39		1,2,4,5-Tetrachlorobenzene	40.000	35.512	11.2	107	-0.09
40	P	Hexachlorocyclopentadiene	40.000	12.597	68.5#	22	-0.09
41	S	2,4,6-Tribromophenol	80.000	87.456	-9.3	127	-0.10
42	C	2,4,6-Trichlorophenol	40.000	42.934	-7.3	127	-0.08
43		2,4,5-Trichlorophenol	40.000	41.682	-4.2	124	-0.07
44	S	2-Fluorobiphenyl	80.000	74.282	7.1	115	-0.08

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF041415\
Data File : BF078527.D
Acq On : 15 Apr 2015 6:18
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 15 05:57:13 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 15 01:10:59 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	37.691	5.8	114	-0.09
46		2-Chloronaphthalene	40.000	37.234	6.9	114	-0.09
47		2-Nitroaniline	40.000	46.813	-17.0	136	-0.08
48		Acenaphthylene	40.000	39.075	2.3	118	-0.10
49		Dimethylphthalate	40.000	39.471	1.3	123	-0.07
50		2,6-Dinitrotoluene	40.000	39.706	0.7	117	-0.08
51	C	Acenaphthene	40.000	38.472	3.8	120	-0.09
52		3-Nitroaniline	40.000	49.944	-24.9	140	-0.08
53	P	2,4-Dinitrophenol	40.000	9.393	76.5#	5	-0.06
54		Dibenzofuran	40.000	40.236	-0.6	125	-0.09
55	P	4-Nitrophenol	40.000	46.545	-16.4	144	-0.05
56		2,4-Dinitrotoluene	40.000	43.249	-8.1	124	-0.08
57		Fluorene	40.000	42.909	-7.3	130	-0.09
58		2,3,4,6-Tetrachlorophenol	40.000	43.295	-8.2	126	-0.09
59		Diethylphthalate	40.000	40.818	-2.0	122	-0.08
60		4-Chlorophenyl-phenylether	40.000	39.593	1.0	122	-0.10
61		4-Nitroaniline	40.000	53.504	-33.8#	148	-0.08
62		Azobenzene	40.000	43.783	-9.5	134	-0.09
63	I	Phenanthrene-d10	20.000	20.000	0.0	140	-0.10
64		4,6-Dinitro-2-methylphenol	40.000	11.077	72.3#	12	-0.07
65	c	n-Nitrosodiphenylamine	40.000	38.580	3.6	132	-0.09
66		4-Bromophenyl-phenylether	40.000	37.272	6.8	127	-0.09
67		Hexachlorobenzene	40.000	34.425	13.9	118	-0.10
68		Atrazine	40.000	37.171	7.1	120	-0.08
69	C	Pentachlorophenol	40.000	55.567	-38.9#	204	-0.09
70		Phenanthrene	40.000	38.074	4.8	129	-0.10
71		Anthracene	40.000	38.635	3.4	131	-0.10
72		Carbazole	40.000	38.320	4.2	126	-0.10
73		Di-n-butylphthalate	40.000	39.084	2.3	127	-0.09
74	C	Fluoranthene	40.000	42.498	-6.2	145	-0.10
75	I	Chrysene-d12	20.000	20.000	0.0	148	-0.11
76		Benzidine	40.000	54.292	-35.7#	196	-0.09
77		Pyrene	40.000	37.230	6.9	138	-0.11
78	S	Terphenyl-d14	80.000	70.405	12.0	129	-0.10
79		Butylbenzylphthalate	40.000	41.667	-4.2	142	-0.10
80		Benzo(a)anthracene	40.000	38.359	4.1	135	-0.11
81		3,3'-Dichlorobenzidine	40.000	43.103	-7.8	153	-0.11
82		Chrysene	40.000	37.960	5.1	145	-0.11
83		Bis(2-ethylhexyl)phthalate	40.000	37.581	6.0	129	-0.11
84	c	Di-n-octyl phthalate	40.000	41.710	-4.3	143	-0.11
85		Indeno(1,2,3-cd)pyrene	40.000	40.070	-0.2	145	-0.21
86	I	Perylene-d12	20.000	20.000	0.0	160	-0.13
87		Benzo(b)fluoranthene	40.000	41.564	-3.9	172	-0.13

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF041415\
Data File : BF078527.D
Acq On : 15 Apr 2015 6:18
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 15 05:57:13 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF040115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 15 01:10:59 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	31.556	21.1	124	-0.12
89 C	Benzo(a)pyrene	40.000	39.996	0.0	157	-0.13
90	Dibenzo(a,h)anthracene	40.000	37.048	7.4	146	-0.23
91	Benzo(g,h,i)perylene	40.000	36.580	8.6	146	-0.26

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