Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071415\
Data File : BF080492.D
Acq On : 15 Jul 2015 14:57
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 15 16:17:47 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 15 03:58:42 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	143	0.00
2		1,4-Dioxane	40.000	41.538	-3.8	159	0.00
3		Pyridine	40.000	46.125	-15.3	162	-0.03
4		n-Nitrosodimethylamine	40.000	46.742	-16.9	165	-0.01
5	S	2-Fluorophenol	80.000	80.158	-0.2	145	0.00
6		Aniline	40.000	39.923	0.2	139	0.00
7	S	Phenol-d6	80.000	82.829	-3.5	149	-0.01
8		2-Chlorophenol	40.000	39.578	1.1	142	0.00
9		Benzaldehyde	40.000	41.666	-4.2	145	0.00
10	C	Phenol	40.000	38.560	3.6	138	0.00
11		bis(2-Chloroethyl)ether	40.000	38.735	3.2	136	0.00
12		1,3-Dichlorobenzene	40.000	40.856	-2.1	147	0.00
13	C	1,4-Dichlorobenzene	40.000	40.821	-2.1	151	0.00
14		1,2-Dichlorobenzene	40.000	39.239	1.9	144	0.00
15		Benzyl Alcohol	40.000	39.748	0.6	139	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.059	-2.6	149	0.00
17		2-Methylphenol	40.000	40.587	-1.5	146	0.00
18		Hexachloroethane	40.000	40.720	-1.8	147	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.752	0.6	135	0.00
20		3+4-Methylphenols	40.000	39.999	0.0	137	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	145	0.00
22		Acetophenone	40.000	40.266	-0.7	151	-0.01
23	S	Nitrobenzene-d5	80.000	79.841	0.2	146	0.00
24		Nitrobenzene	40.000	38.399	4.0	140	0.00
25		Isophorone	40.000	41.565	-3.9	154	0.00
26	C	2-Nitrophenol	40.000	39.392	1.5	140	0.00
27		2,4-Dimethylphenol	40.000	40.112	-0.3	142	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.687	0.8	152	0.00
29	C	2,4-Dichlorophenol	40.000	40.154	-0.4	154	0.00
30		1,2,4-Trichlorobenzene	40.000	38.366	4.1	143	0.00
31		Naphthalene	40.000	38.528	3.7	146	0.00
32		Benzoic acid	40.000	45.890	-14.7	167	0.00
33		4-Chloroaniline	40.000	40.074	-0.2	143	0.00
34	C	Hexachlorobutadiene	40.000	38.464	3.8	138	0.00
35		Caprolactam	40.000	42.176	-5.4	156	-0.01
36	C	4-Chloro-3-methylphenol	40.000	40.438	-1.1	153	-0.01
37		2-Methylnaphthalene	40.000	36.563	8.6	133	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	139	-0.01
39		1,2,4,5-Tetrachlorobenzene	40.000	38.634	3.4	147	0.00
40	P	Hexachlorocyclopentadiene	40.000	28.944	27.6#	92	0.00
41	S	2,4,6-Tribromophenol	80.000	80.494	-0.6	134	0.00
42	C	2,4,6-Trichlorophenol	40.000	39.312	1.7	133	0.00
43		2,4,5-Trichlorophenol	40.000	40.136	-0.3	139	-0.01
44	S	2-Fluorobiphenyl	80.000	77.062	3.7	143	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071415\
Data File : BF080492.D
Acq On : 15 Jul 2015 14:57
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 15 16:17:47 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 15 03:58:42 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	39.023	2.4	148	0.00
46		2-Chloronaphthalene	40.000	38.200	4.5	134	0.00
47		2-Nitroaniline	40.000	41.449	-3.6	149	0.00
48		Acenaphthylene	40.000	39.528	1.2	143	0.00
49		Dimethylphthalate	40.000	40.412	-1.0	150	0.00
50		2,6-Dinitrotoluene	40.000	41.104	-2.8	148	0.00
51	C	Acenaphthene	40.000	38.192	4.5	137	0.00
52		3-Nitroaniline	40.000	41.573	-3.9	150	0.00
53	P	2,4-Dinitrophenol	40.000	26.918	32.7#	76	0.00
54		Dibenzofuran	40.000	38.133	4.7	142	0.00
55	P	4-Nitrophenol	40.000	39.308	1.7	137	0.00
56		2,4-Dinitrotoluene	40.000	40.843	-2.1	149	-0.01
57		Fluorene	40.000	39.579	1.1	141	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	43.306	-8.3	157	0.00
59		Diethylphthalate	40.000	40.169	-0.4	149	0.00
60		4-Chlorophenyl-phenylether	40.000	40.244	-0.6	150	0.00
61		4-Nitroaniline	40.000	43.893	-9.7	152	0.00
62		Azobenzene	40.000	41.567	-3.9	149	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	141	0.00
64		4,6-Dinitro-2-methylphenol	40.000	26.053	34.9#	86	0.00
65	c	n-Nitrosodiphenylamine	40.000	37.501	6.2	137	0.00
66		4-Bromophenyl-phenylether	40.000	39.116	2.2	151	0.00
67		Hexachlorobenzene	40.000	36.326	9.2	131	0.00
68		Atrazine	40.000	42.310	-5.8	159	0.00
69	C	Pentachlorophenol	40.000	41.210	-3.0	149	0.00
70		Phenanthrene	40.000	36.377	9.1	137	0.00
71		Anthracene	40.000	41.361	-3.4	151	0.00
72		Carbazole	40.000	41.006	-2.5	154	0.00
73		Di-n-butylphthalate	40.000	46.227	-15.6	178	0.02
74	C	Fluoranthene	40.000	41.950	-4.9	158	0.05
75	I	Chrysene-d12	20.000	20.000	0.0	157	0.14
76		Benzidine	40.000	37.182	7.0	138	0.06
77		Pyrene	40.000	39.620	1.0	158	0.06
78	S	Terphenyl-d14	80.000	80.397	-0.5	157	0.07
79		Butylbenzylphthalate	40.000	44.062	-10.2	171	0.10
80		Benzo(a)anthracene	40.000	39.519	1.2	167	0.14
81		3,3'-Dichlorobenzidine	40.000	40.933	-2.3	163	0.14
82		Chrysene	40.000	38.161	4.6	151	0.14
83		Bis(2-ethylhexyl)phthalate	40.000	45.551	-13.9	179	0.14
84	c	Di-n-octyl phthalate	40.000	43.194	-8.0	170	0.17
85		Indeno(1,2,3-cd)pyrene	40.000	30.267	24.3	111	0.22
86	I	Perylene-d12	20.000	20.000	0.0	140	0.21
87		Benzo(b)fluoranthene	40.000	39.168	2.1	152	0.19

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071415\
Data File : BF080492.D
Acq On : 15 Jul 2015 14:57
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 15 16:17:47 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 15 03:58:42 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	40.435	-1.1	138	0.19
89 C	Benzo(a)pyrene	40.000	37.374	6.6	135	0.19
90	Dibenzo(a,h)anthracene	40.000	32.231	19.4	113	0.22
91	Benzo(g,h,i)perylene	40.000	29.691	25.8#	104	0.21

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