Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF122817\
Data File : BF101795.D
Acq On : 28 Dec 2017 11:51
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 28 12:50:12 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 26 10:48:01 2017
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	87	0.00
2		1,4-Dioxane	40.000	35.633	10.9	77	-0.01
3		Pyridine	40.000	34.672	13.3	74	-0.02
4		n-Nitrosodimethylamine	40.000	35.234	11.9	75	-0.01
5	S	2-Fluorophenol	80.000	80.724	-0.9	88	0.00
6		Aniline	40.000	35.921	10.2	80	0.00
7	S	Phenol-d6	80.000	73.643	7.9	84	0.00
8		2-Chlorophenol	40.000	38.833	2.9	87	0.00
9		Benzaldehyde	40.000	33.190	17.0	73	0.00
10	C	Phenol	40.000	36.006	10.0	80	0.00
11		bis(2-Chloroethyl)ether	40.000	37.528	6.2	83	0.00
12		1,3-Dichlorobenzene	40.000	39.231	1.9	87	0.00
13	C	1,4-Dichlorobenzene	40.000	39.477	1.3	89	0.00
14		1,2-Dichlorobenzene	40.000	38.758	3.1	86	0.00
15		Benzyl Alcohol	40.000	37.492	6.3	84	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	37.232	6.9	81	0.00
17		2-Methylphenol	40.000	36.864	7.8	85	0.00
18		Hexachloroethane	40.000	38.866	2.8	86	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	37.114	7.2	87	0.00
20		3+4-Methylphenols	40.000	43.553	-8.9	96	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	87	0.00
22		Acetophenone	40.000	39.523	1.2	88	0.00
23	S	Nitrobenzene-d5	80.000	76.878	3.9	84	0.00
24		Nitrobenzene	40.000	37.221	6.9	83	0.00
25		Isophorone	40.000	36.575	8.6	82	0.00
26	C	2-Nitrophenol	40.000	44.708	-11.8	95	0.00
27		2,4-Dimethylphenol	40.000	38.681	3.3	87	0.00
28		bis(2-Chloroethoxy)methane	40.000	37.529	6.2	85	0.00
29	C	2,4-Dichlorophenol	40.000	41.464	-3.7	91	0.00
30		1,2,4-Trichlorobenzene	40.000	38.484	3.8	85	0.00
31		Naphthalene	40.000	38.061	4.8	86	0.00
32		Benzoic acid	40.000	46.207	-15.5	117	0.00
33		4-Chloroaniline	40.000	38.795	3.0	88	0.00
34	C	Hexachlorobutadiene	40.000	40.316	-0.8	90	0.00
35		Caprolactam	40.000	39.697	0.8	88	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.379	1.6	87	0.00
37		2-Methylnaphthalene	40.000	37.997	5.0	87	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	90	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	39.141	2.1	90	0.00
40	P	Hexachlorocyclopentadiene	40.000	33.289	16.8	70	0.00
41	S	2,4,6-Tribromophenol	80.000	86.021	-7.5	97	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.412	-1.0	90	0.00
43		2,4,5-Trichlorophenol	40.000	36.833	7.9	82	0.00
44	S	2-Fluorobiphenyl	80.000	76.556	4.3	88	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF122817\
Data File : BF101795.D
Acq On : 28 Dec 2017 11:51
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 28 12:50:12 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 26 10:48:01 2017
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	37.151	7.1	88	0.00
46		2-Chloronaphthalene	40.000	37.189	7.0	88	0.00
47		2-Nitroaniline	40.000	39.179	2.1	88	0.00
48		Acenaphthylene	40.000	36.576	8.6	87	0.00
49		Dimethylphthalate	40.000	36.304	9.2	86	0.00
50		2,6-Dinitrotoluene	40.000	39.313	1.7	87	0.00
51	C	Acenaphthene	40.000	37.453	6.4	89	0.00
52		3-Nitroaniline	40.000	40.862	-2.2	91	0.00
53	P	2,4-Dinitrophenol	40.000	44.075	-10.2	112	0.00
54		Dibenzofuran	40.000	41.309	-3.3	94	0.00
55	P	4-Nitrophenol	40.000	40.877	-2.2	88	0.00
56		2,4-Dinitrotoluene	40.000	48.403	-21.0	108	0.00
57		Fluorene	40.000	39.977	0.1	91	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	41.249	-3.1	92	0.00
59		Diethylphthalate	40.000	38.585	3.5	93	0.00
60		4-Chlorophenyl-phenylether	40.000	41.200	-3.0	92	0.00
61		4-Nitroaniline	40.000	40.026	-0.1	91	0.00
62		Azobenzene	40.000	35.877	10.3	86	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	94	0.00
64		4,6-Dinitro-2-methylphenol	40.000	47.525	-18.8	106	0.00
65	c	n-Nitrosodiphenylamine	40.000	36.891	7.8	90	0.00
66		4-Bromophenyl-phenylether	40.000	39.087	2.3	94	0.00
67		Hexachlorobenzene	40.000	38.312	4.2	92	0.00
68		Atrazine	40.000	38.312	4.2	91	0.00
69	C	Pentachlorophenol	40.000	35.649	10.9	87	0.00
70		Phenanthrene	40.000	36.965	7.6	92	0.00
71		Anthracene	40.000	37.452	6.4	93	0.00
72		Carbazole	40.000	37.960	5.1	93	0.00
73		Di-n-butylphthalate	40.000	39.116	2.2	97	0.00
74	C	Fluoranthene	40.000	38.201	4.5	95	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	96	0.00
76		Benzidine	40.000	44.490	-11.2	108	0.00
77		Pyrene	40.000	37.489	6.3	94	0.00
78	S	Terphenyl-d14	80.000	74.252	7.2	96	0.00
79		Butylbenzylphthalate	40.000	39.534	1.2	96	0.00
80		Benzo(a)anthracene	40.000	36.891	7.8	92	0.00
81		3,3'-Dichlorobenzidine	40.000	36.642	8.4	89	0.00
82		Chrysene	40.000	37.574	6.1	94	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	37.339	6.7	92	0.00
84	c	Di-n-octyl phthalate	40.000	38.910	2.7	96	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	37.906	5.2	97	0.01
86	I	Perylene-d12	20.000	20.000	0.0	106	0.00
87		Benzo(b)fluoranthene	40.000	33.693	15.8	89	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF122817\
Data File : BF101795.D
Acq On : 28 Dec 2017 11:51
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Dec 28 12:50:12 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF121417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Dec 26 10:48:01 2017
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	35.198	12.0	99	0.00
89 C	Benzo(a)pyrene	40.000	34.329	14.2	93	0.00
90	Dibenzo(a,h)anthracene	40.000	34.581	13.5	97	0.01
91	Benzo(g,h,i)perylene	40.000	35.738	10.7	99	0.01

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