Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042619\
Data File : BF114015.D
Acq On : 26 Apr 2019 17:53
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 26 18:16:02 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 23 03:32:35 2019
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	83	0.00
2		1,4-Dioxane	40.000	40.151	-0.4	83	0.01
3		Pyridine	40.000	39.133	2.2	81	0.00
4		n-Nitrosodimethylamine	40.000	33.058	17.4	69	0.00
5	S	2-Fluorophenol	80.000	80.639	-0.8	82	0.00
6		Aniline	40.000	42.171	-5.4	87	0.00
7	S	Phenol-d6	80.000	85.208	-6.5	87	0.00
8		2-Chlorophenol	40.000	42.732	-6.8	87	0.00
9		Benzaldehyde	40.000	42.909	-7.3	86	0.00
10	C	Phenol	40.000	42.125	-5.3	86	0.00
11		bis(2-Chloroethyl)ether	40.000	42.263	-5.7	87	0.00
12		1,3-Dichlorobenzene	40.000	42.827	-7.1	88	0.00
13	C	1,4-Dichlorobenzene	40.000	42.505	-6.3	87	0.00
14		1,2-Dichlorobenzene	40.000	43.036	-7.6	87	0.00
15		Benzyl Alcohol	40.000	48.900	-22.2	99	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	40.302	-0.8	83	0.00
17		2-Methylphenol	40.000	41.410	-3.5	85	0.00
18		Hexachloroethane	40.000	43.242	-8.1	89	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	42.403	-6.0	88	0.00
20		3+4-Methylphenols	40.000	43.292	-8.2	87	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	83	0.00
22		Acetophenone	40.000	42.533	-6.3	88	0.00
23	S	Nitrobenzene-d5	80.000	90.328	-12.9	92	0.00
24		Nitrobenzene	40.000	43.914	-9.8	89	0.00
25		Isophorone	40.000	42.057	-5.1	89	0.00
26	C	2-Nitrophenol	40.000	49.906	-24.8#	102	0.00
27		2,4-Dimethylphenol	40.000	40.803	-2.0	83	0.00
28		bis(2-Chloroethoxy)methane	40.000	42.567	-6.4	88	0.00
29	C	2,4-Dichlorophenol	40.000	43.219	-8.0	89	0.00
30		1,2,4-Trichlorobenzene	40.000	42.750	-6.9	88	0.00
31		Naphthalene	40.000	42.736	-6.8	87	0.00
32		Benzoic acid	40.000	34.458	13.9	68	0.02
33		4-Chloroaniline	40.000	41.959	-4.9	88	0.00
34	C	Hexachlorobutadiene	40.000	42.992	-7.5	88	0.00
35		Caprolactam	40.000	42.774	-6.9	91	0.02
36	C	4-Chloro-3-methylphenol	40.000	43.634	-9.1	90	0.00
37		2-Methylnaphthalene	40.000	43.072	-7.7	88	0.00
38		1-Methylnaphthalene	40.000	42.944	-7.4	89	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	85	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	42.334	-5.8	89	0.00
41	P	Hexachlorocyclopentadiene	40.000	38.021	4.9	79	0.00
42	S	2,4,6-Tribromophenol	80.000	86.610	-8.3	90	0.00
43	C	2,4,6-Trichlorophenol	40.000	42.906	-7.3	91	0.00
44		2,4,5-Trichlorophenol	40.000	42.727	-6.8	90	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042619\
Data File : BF114015.D
Acq On : 26 Apr 2019 17:53
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 26 18:16:02 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 23 03:32:35 2019
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	S	2-Fluorobiphenyl	80.000	83.089	-3.9	90	0.00
46		1,1'-Biphenyl	40.000	42.837	-7.1	90	0.00
47		2-Chloronaphthalene	40.000	42.574	-6.4	90	0.00
48		2-Nitroaniline	40.000	46.281	-15.7	96	0.00
49		Acenaphthylene	40.000	42.866	-7.2	90	0.00
50		Dimethylphthalate	40.000	44.160	-10.4	95	0.00
51		2,6-Dinitrotoluene	40.000	47.081	-17.7	97	0.00
52	C	Acenaphthene	40.000	43.946	-9.9	93	0.00
53		3-Nitroaniline	40.000	44.328	-10.8	93	0.00
54	P	2,4-Dinitrophenol	40.000	52.200	-30.5#	130	0.00
55		Dibenzofuran	40.000	43.105	-7.8	90	0.00
56	P	4-Nitrophenol	40.000	46.069	-15.2	96	0.00
57		2,4-Dinitrotoluene	40.000	50.553	-26.4#	106	0.00
58		Fluorene	40.000	43.211	-8.0	91	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	43.589	-9.0	92	0.00
60		Diethylphthalate	40.000	44.261	-10.7	93	0.00
61		4-Chlorophenyl-phenylether	40.000	43.197	-8.0	90	0.00
62		4-Nitroaniline	40.000	47.666	-19.2	101	0.00
63		Azobenzene	40.000	42.764	-6.9	90	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	89	0.00
65		4,6-Dinitro-2-methylphenol	40.000	50.749	-26.9#	126	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.830	-4.6	93	0.00
67		4-Bromophenyl-phenylether	40.000	41.262	-3.2	90	0.00
68		Hexachlorobenzene	40.000	41.669	-4.2	90	0.00
69		Atrazine	40.000	40.112	-0.3	89	0.00
70	C	Pentachlorophenol	40.000	43.433	-8.6	94	0.00
71		Phenanthrene	40.000	41.938	-4.8	91	0.00
72		Anthracene	40.000	42.224	-5.6	92	0.00
73		Carbazole	40.000	43.213	-8.0	94	0.00
74		Di-n-butylphthalate	40.000	44.298	-10.7	96	0.00
75	C	Fluoranthene	40.000	41.120	-2.8	93	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	83	0.00
77		Benzidine	40.000	43.320	-8.3	96	0.00
78		Pyrene	40.000	46.134	-15.3	93	0.00
79	S	Terphenyl-d14	80.000	92.537	-15.7	93	0.00
80		Butylbenzylphthalate	40.000	50.492	-26.2#	103	0.00
81		Benzo(a)anthracene	40.000	47.368	-18.4	97	0.00
82		3,3'-Dichlorobenzidine	40.000	48.094	-20.2	96	0.00
83		Chrysene	40.000	46.635	-16.6	95	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	52.333	-30.8#	108	0.00
85	c	Di-n-octyl phthalate	40.000	51.672	-29.2#	107	-0.03
86		Indeno(1,2,3-cd)pyrene	40.000	37.894	5.3	84	-0.02
87	I	Perylene-d12	20.000	20.000	0.0	84	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF042619\
Data File : BF114015.D
Acq On : 26 Apr 2019 17:53
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 26 18:16:02 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF042219.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Apr 23 03:32:35 2019
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(b)fluoranthene	40.000	46.539	-16.3	98	-0.02
89	Benzo(k)fluoranthene	40.000	43.559	-8.9	89	-0.02
90 C	Benzo(a)pyrene	40.000	43.524	-8.8	91	-0.02
91	Dibenzo(a,h)anthracene	40.000	39.556	1.1	84	-0.02
92	Benzo(g,h,i)perylene	40.000	38.696	3.3	83	-0.02

(#) = Out of Range SPCC's out = 0 CCC's out = 2