Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040919\
Data File : BN005114.D
Acq On : 09 Apr 2019 16:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Apr 10 01:15:50 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN032619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 27 05:27:09 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	89	-0.02
2		1,4-Dioxane	40.000	36.652	8.4	82	-0.03
3		Pyridine	40.000	39.612	1.0	90	-0.03
4		n-Nitrosodimethylamine	40.000	36.608	8.5	80	-0.03
5	S	2-Fluorophenol	80.000	79.051	1.2	89	-0.03
6		Aniline	40.000	40.418	-1.0	89	-0.02
7	S	Phenol-d6	80.000	81.068	-1.3	89	-0.03
8		2-Chlorophenol	40.000	41.064	-2.7	91	-0.03
9		Benzaldehyde	40.000	38.372	4.1	82	-0.03
10	C	Phenol	40.000	40.133	-0.3	88	-0.02
11		bis(2-Chloroethyl)ether	40.000	40.391	-1.0	89	-0.03
12		1,3-Dichlorobenzene	40.000	40.315	-0.8	90	-0.03
13	C	1,4-Dichlorobenzene	40.000	40.470	-1.2	90	-0.03
14		1,2-Dichlorobenzene	40.000	41.397	-3.5	91	-0.03
15		Benzyl Alcohol	40.000	40.255	-0.6	88	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	37.338	6.7	83	-0.02
17		2-Methylphenol	40.000	41.383	-3.5	90	-0.03
18		Hexachloroethane	40.000	39.695	0.8	88	-0.03
19	P	n-Nitroso-di-n-propylamine	40.000	40.363	-0.9	87	-0.02
20		3+4-Methylphenols	40.000	41.181	-3.0	90	-0.03
21	I	Naphthalene-d8	20.000	20.000	0.0	89	-0.03
22		Acetophenone	40.000	40.501	-1.3	90	-0.03
23	S	Nitrobenzene-d5	80.000	78.334	2.1	88	-0.03
24		Nitrobenzene	40.000	39.000	2.5	87	-0.03
25		Isophorone	40.000	39.194	2.0	87	-0.02
26	C	2-Nitrophenol	40.000	40.787	-2.0	91	-0.02
27		2,4-Dimethylphenol	40.000	40.518	-1.3	90	-0.03
28		bis(2-Chloroethoxy)methane	40.000	40.446	-1.1	90	-0.02
29	C	2,4-Dichlorophenol	40.000	41.903	-4.8	93	-0.03
30		1,2,4-Trichlorobenzene	40.000	41.635	-4.1	94	-0.03
31		Naphthalene	40.000	41.043	-2.6	92	-0.03
32		Benzoic acid	40.000	33.980	15.1	72	-0.02
33		4-Chloroaniline	40.000	41.751	-4.4	92	-0.03
34	C	Hexachlorobutadiene	40.000	41.331	-3.3	93	-0.03
35		Caprolactam	40.000	41.060	-2.7	87	-0.02
36	C	4-Chloro-3-methylphenol	40.000	40.415	-1.0	88	-0.02
37		2-Methylnaphthalene	40.000	41.890	-4.7	93	-0.02
38		1-Methylnaphthalene	40.000	41.669	-4.2	92	-0.02
39	I	Acenaphthene-d10	20.000	20.000	0.0	91	-0.02
40		1,2,4,5-Tetrachlorobenzene	40.000	41.078	-2.7	94	-0.02
41	P	Hexachlorocyclopentadiene	40.000	42.640	-6.6	100	-0.02
42	S	2,4,6-Tribromophenol	80.000	88.913	-11.1	99	-0.02
43	C	2,4,6-Trichlorophenol	40.000	40.394	-1.0	92	-0.02
44		2,4,5-Trichlorophenol	40.000	39.946	0.1	89	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040919\
Data File : BN005114.D
Acq On : 09 Apr 2019 16:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Apr 10 01:15:50 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN032619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 27 05:27:09 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	82.818	-3.5	94	-0.02
46		1,1'-Biphenyl	40.000	41.240	-3.1	94	-0.02
47		2-Chloronaphthalene	40.000	41.022	-2.6	94	-0.03
48		2-Nitroaniline	40.000	37.886	5.3	85	-0.02
49		Acenaphthylene	40.000	40.547	-1.4	92	-0.03
50		Dimethylphthalate	40.000	41.350	-3.4	93	-0.02
51		2,6-Dinitrotoluene	40.000	40.796	-2.0	92	-0.02
52	C	Acenaphthene	40.000	41.188	-3.0	93	-0.02
53		3-Nitroaniline	40.000	39.575	1.1	88	-0.02
54	P	2,4-Dinitrophenol	40.000	41.025	-2.6	92	-0.02
55		Dibenzofuran	40.000	40.974	-2.4	93	-0.02
56	P	4-Nitrophenol	40.000	34.720	13.2	78	-0.02
57		2,4-Dinitrotoluene	40.000	41.393	-3.5	92	-0.02
58		Fluorene	40.000	41.134	-2.8	92	-0.02
59		2,3,4,6-Tetrachlorophenol	40.000	41.094	-2.7	92	-0.02
60		Diethylphthalate	40.000	41.252	-3.1	92	-0.02
61		4-Chlorophenyl-phenylether	40.000	41.604	-4.0	94	-0.02
62		4-Nitroaniline	40.000	39.166	2.1	86	-0.02
63		Azobenzene	40.000	39.030	2.4	88	-0.02
64	I	Phenanthrene-d10	20.000	20.000	0.0	90	-0.02
65		4,6-Dinitro-2-methylphenol	40.000	37.280	6.8	83	-0.02
66	c	n-Nitrosodiphenylamine	40.000	40.470	-1.2	93	-0.02
67		4-Bromophenyl-phenylether	40.000	41.722	-4.3	95	-0.02
68		Hexachlorobenzene	40.000	43.104	-7.8	98	-0.02
69		Atrazine	40.000	40.343	-0.9	90	-0.02
70	C	Pentachlorophenol	40.000	40.009	-0.0	93	-0.02
71		Phenanthrene	40.000	40.618	-1.5	92	-0.02
72		Anthracene	40.000	41.306	-3.3	94	-0.02
73		Carbazole	40.000	40.613	-1.5	91	-0.02
74		Di-n-butylphthalate	40.000	42.287	-5.7	95	-0.02
75	C	Fluoranthene	40.000	41.904	-4.8	94	-0.02
76	I	Chrysene-d12	20.000	20.000	0.0	97	-0.02
77		Benzidine	40.000	38.197	4.5	83	-0.02
78		Pyrene	40.000	39.905	0.2	93	-0.02
79	S	Terphenyl-d14	80.000	85.779	-7.2	100	-0.02
80		Butylbenzylphthalate	40.000	42.617	-6.5	101	-0.02
81		Benzo(a)anthracene	40.000	40.513	-1.3	97	-0.02
82		3,3'-Dichlorobenzidine	40.000	40.070	-0.2	95	-0.02
83		Chrysene	40.000	40.203	-0.5	97	-0.02
84		Bis(2-ethylhexyl)phthalate	40.000	45.667	-14.2	108	-0.02
85	c	Di-n-octyl phthalate	40.000	45.964	-14.9	112	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	35.785	10.5	94	-0.04
87	I	Perylene-d12	20.000	20.000	0.0	94	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN040919\
Data File : BN005114.D
Acq On : 09 Apr 2019 16:04
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Apr 10 01:15:50 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\8270-BN032619.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Mar 27 05:27:09 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	42.373	-5.9	100	-0.02
89	Benzo(k)fluoranthene	40.000	41.536	-3.8	95	-0.02
90 C	Benzo(a)pyrene	40.000	40.588	-1.5	95	-0.03
91	Dibenzo(a,h)anthracene	40.000	39.293	1.8	95	-0.04
92	Benzo(g,h,i)perylene	40.000	38.540	3.7	94	-0.05

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