Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102418\
Data File : BN003288.D
Acq On : 25 Oct 2018 00:59
Operator : MJ/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Oct 25 05:26:26 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN102418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 24 14:21:25 2018
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	111	0.00
2		1,4-Dioxane	40.000	37.664	5.8	107	0.00
3		Pyridine	40.000	40.943	-2.4	107	0.00
4		n-Nitrosodimethylamine	40.000	38.624	3.4	108	0.00
5	S	2-Fluorophenol	80.000	78.185	2.3	107	0.00
6		Aniline	40.000	40.509	-1.3	110	0.00
7	S	Phenol-d6	80.000	80.145	-0.2	109	0.00
8		2-Chlorophenol	40.000	39.495	1.3	109	0.00
9		Benzaldehyde	40.000	38.792	3.0	107	0.00
10	C	Phenol	40.000	39.901	0.2	108	0.00
11		bis(2-Chloroethyl)ether	40.000	39.337	1.7	108	0.00
12		1,3-Dichlorobenzene	40.000	39.245	1.9	108	0.00
13	C	1,4-Dichlorobenzene	40.000	39.305	1.7	109	0.00
14		1,2-Dichlorobenzene	40.000	38.812	3.0	107	0.00
15		Benzyl Alcohol	40.000	40.380	-1.0	113	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	39.333	1.7	110	0.00
17		2-Methylphenol	40.000	40.008	-0.0	109	0.00
18		Hexachloroethane	40.000	39.028	2.4	109	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.622	-1.6	111	0.00
20		3+4-Methylphenols	40.000	40.210	-0.5	109	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	113	0.00
22		Acetophenone	40.000	39.950	0.1	111	0.00
23	S	Nitrobenzene-d5	80.000	79.182	1.0	110	0.00
24		Nitrobenzene	40.000	39.535	1.2	110	0.00
25		Isophorone	40.000	40.282	-0.7	112	0.00
26	C	2-Nitrophenol	40.000	40.444	-1.1	110	0.00
27		2,4-Dimethylphenol	40.000	39.314	1.7	109	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.922	0.2	112	0.00
29	C	2,4-Dichlorophenol	40.000	40.622	-1.6	111	0.00
30		1,2,4-Trichlorobenzene	40.000	39.108	2.2	110	0.00
31		Naphthalene	40.000	38.933	2.7	110	0.00
32		Benzoic acid	40.000	36.357	9.1	104	0.00
33		4-Chloroaniline	40.000	39.707	0.7	111	0.00
34	C	Hexachlorobutadiene	40.000	39.265	1.8	110	0.00
35		Caprolactam	40.000	40.997	-2.5	109	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.396	-1.0	111	0.00
37		2-Methylnaphthalene	40.000	39.450	1.4	111	0.00
38		1-Methylnaphthalene	40.000	39.279	1.8	111	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	116	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	39.275	1.8	111	0.00
41	P	Hexachlorocyclopentadiene	40.000	38.847	2.9	117	0.00
42	S	2,4,6-Tribromophenol	80.000	80.226	-0.3	113	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.015	-0.0	112	0.00
44		2,4,5-Trichlorophenol	40.000	40.182	-0.5	115	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102418\
Data File : BN003288.D
Acq On : 25 Oct 2018 00:59
Operator : MJ/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Oct 25 05:26:26 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN102418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 24 14:21:25 2018
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	77.346	3.3	111	0.00
46		1,1'-Biphenyl	40.000	38.847	2.9	112	0.00
47		2-Chloronaphthalene	40.000	39.043	2.4	111	0.00
48		2-Nitroaniline	40.000	40.488	-1.2	111	0.00
49		Acenaphthylene	40.000	39.210	2.0	112	0.00
50		Dimethylphthalate	40.000	38.738	3.2	112	0.00
51		2,6-Dinitrotoluene	40.000	39.408	1.5	112	0.00
52	C	Acenaphthene	40.000	38.980	2.6	112	0.00
53		3-Nitroaniline	40.000	40.041	-0.1	111	0.00
54	P	2,4-Dinitrophenol	40.000	36.026	9.9	105	0.00
55		Dibenzofuran	40.000	38.342	4.1	111	0.00
56	P	4-Nitrophenol	40.000	37.662	5.8	109	0.00
57		2,4-Dinitrotoluene	40.000	40.185	-0.5	112	0.00
58		Fluorene	40.000	38.619	3.5	111	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.226	-0.6	114	0.00
60		Diethylphthalate	40.000	39.295	1.8	113	0.00
61		4-Chlorophenyl-phenylether	40.000	38.998	2.5	113	0.00
62		4-Nitroaniline	40.000	40.987	-2.5	110	0.00
63		Azobenzene	40.000	39.283	1.8	113	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	117	0.00
65		4,6-Dinitro-2-methylphenol	40.000	38.778	3.1	111	0.00
66	c	n-Nitrosodiphenylamine	40.000	38.520	3.7	112	0.00
67		4-Bromophenyl-phenylether	40.000	39.154	2.1	112	0.00
68		Hexachlorobenzene	40.000	38.660	3.4	111	0.00
69		Atrazine	40.000	39.321	1.7	111	0.00
70	C	Pentachlorophenol	40.000	41.376	-3.4	125	0.00
71		Phenanthrene	40.000	38.383	4.0	111	0.00
72		Anthracene	40.000	38.721	3.2	111	0.00
73		Carbazole	40.000	38.535	3.7	111	0.00
74		Di-n-butylphthalate	40.000	38.654	3.4	111	0.00
75	C	Fluoranthene	40.000	38.511	3.7	112	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	117	0.00
77		Benzidine	40.000	42.536	-6.3	123	0.00
78		Pyrene	40.000	38.466	3.8	112	0.00
79	S	Terphenyl-d14	80.000	76.299	4.6	110	0.00
80		Butylbenzylphthalate	40.000	38.434	3.9	110	0.00
81		Benzo(a)anthracene	40.000	37.754	5.6	109	0.00
82		3,3'-Dichlorobenzidine	40.000	37.720	5.7	108	0.00
83		Chrysene	40.000	38.222	4.4	111	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	38.207	4.5	111	0.00
85	c	Di-n-octyl phthalate	40.000	37.435	6.4	108	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	35.282	11.8	99	0.00
87	I	Perylene-d12	20.000	20.000	0.0	111	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102418\
Data File : BN003288.D
Acq On : 25 Oct 2018 00:59
Operator : MJ/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Oct 25 05:26:26 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN102418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 24 14:21:25 2018
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	37.889	5.3	105	0.00
89	Benzo(k)fluoranthene	40.000	38.679	3.3	110	0.00
90 C	Benzo(a)pyrene	40.000	38.127	4.7	106	0.00
91	Dibenzo(a,h)anthracene	40.000	36.142	9.6	99	0.00
92	Benzo(g,h,i)perylene	40.000	35.988	10.0	98	0.00

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