Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072618\
Data File : BN002167.D
Acq On : 26 Jul 2018 10:08
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Jul 27 00:56:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN072518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 26 18:27:25 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	106	0.00
2		1,4-Dioxane	40.000	40.055	-0.1	107	0.00
3		Pyridine	40.000	41.985	-5.0	108	0.00
4		n-Nitrosodimethylamine	40.000	39.745	0.6	109	0.00
5	S	2-Fluorophenol	80.000	79.712	0.4	105	0.00
6		Aniline	40.000	40.925	-2.3	109	0.00
7	S	Phenol-d6	80.000	81.482	-1.9	107	0.00
8		2-Chlorophenol	40.000	40.470	-1.2	107	0.00
9		Benzaldehyde	40.000	41.554	-3.9	112	0.00
10	C	Phenol	40.000	41.085	-2.7	109	0.00
11		bis(2-Chloroethyl)ether	40.000	40.594	-1.5	108	0.00
12		1,3-Dichlorobenzene	40.000	39.881	0.3	105	0.00
13	C	1,4-Dichlorobenzene	40.000	39.805	0.5	106	0.00
14		1,2-Dichlorobenzene	40.000	39.471	1.3	106	0.00
15		Benzyl Alcohol	40.000	40.981	-2.5	109	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.473	-3.7	113	0.00
17		2-Methylphenol	40.000	39.896	0.3	106	0.00
18		Hexachloroethane	40.000	39.986	0.0	106	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.853	-2.1	113	0.00
20		3+4-Methylphenols	40.000	40.648	-1.6	109	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	112	0.00
22		Acetophenone	40.000	39.795	0.5	111	0.00
23	S	Nitrobenzene-d5	80.000	83.709	-4.6	115	0.00
24		Nitrobenzene	40.000	39.983	0.0	111	0.00
25		Isophorone	40.000	39.946	0.1	114	0.00
26	C	2-Nitrophenol	40.000	39.702	0.7	109	0.00
27		2,4-Dimethylphenol	40.000	39.610	1.0	109	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.642	0.9	112	0.00
29	C	2,4-Dichlorophenol	40.000	39.916	0.2	111	0.00
30		1,2,4-Trichlorobenzene	40.000	39.285	1.8	109	0.00
31		Naphthalene	40.000	39.169	2.1	110	0.00
32		Benzoic acid	40.000	35.596	11.0	103	0.00
33		4-Chloroaniline	40.000	39.886	0.3	112	0.00
34	C	Hexachlorobutadiene	40.000	39.055	2.4	108	0.00
35		Caprolactam	40.000	39.651	0.9	116	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.150	-0.4	115	0.00
37		2-Methylnaphthalene	40.000	39.664	0.8	113	0.00
38	I	Acenaphthene-d10	20.000	20.000	0.0	117	0.00
39		1,2,4,5-Tetrachlorobenzene	40.000	40.026	-0.1	112	0.00
40	P	Hexachlorocyclopentadiene	40.000	41.456	-3.6	110	0.00
41	S	2,4,6-Tribromophenol	80.000	79.111	1.1	117	0.00
42	C	2,4,6-Trichlorophenol	40.000	40.244	-0.6	114	0.00
43		2,4,5-Trichlorophenol	40.000	40.044	-0.1	114	0.00
44	S	2-Fluorobiphenyl	80.000	82.697	-3.4	118	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072618\
Data File : BN002167.D
Acq On : 26 Jul 2018 10:08
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Jul 27 00:56:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN072518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 26 18:27:25 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		1,1'-Biphenyl	40.000	39.910	0.2	115	0.00
46		2-Chloronaphthalene	40.000	40.103	-0.3	114	0.00
47		2-Nitroaniline	40.000	40.722	-1.8	117	0.00
48		Acenaphthylene	40.000	39.809	0.5	115	0.00
49		Dimethylphthalate	40.000	39.193	2.0	117	0.00
50		2,6-Dinitrotoluene	40.000	40.079	-0.2	117	0.00
51	C	Acenaphthene	40.000	39.927	0.2	116	0.00
52		3-Nitroaniline	40.000	40.112	-0.3	116	0.00
53	P	2,4-Dinitrophenol	40.000	37.366	6.6	112	0.00
54		Dibenzofuran	40.000	39.484	1.3	116	0.00
55	P	4-Nitrophenol	40.000	39.639	0.9	114	0.00
56		2,4-Dinitrotoluene	40.000	40.126	-0.3	118	0.00
57		Fluorene	40.000	39.349	1.6	116	0.00
58		2,3,4,6-Tetrachlorophenol	40.000	39.644	0.9	116	0.00
59		Diethylphthalate	40.000	39.427	1.4	119	0.00
60		4-Chlorophenyl-phenylether	40.000	39.363	1.6	116	0.00
61		4-Nitroaniline	40.000	39.679	0.8	116	0.00
62		Azobenzene	40.000	40.197	-0.5	119	0.00
63	I	Phenanthrene-d10	20.000	20.000	0.0	118	0.00
64		4,6-Dinitro-2-methylphenol	40.000	40.268	-0.7	113	0.00
65	c	n-Nitrosodiphenylamine	40.000	40.345	-0.9	118	0.00
66		4-Bromophenyl-phenylether	40.000	39.994	0.0	116	0.00
67		Hexachlorobenzene	40.000	40.075	-0.2	117	0.00
68		Atrazine	40.000	40.324	-0.8	119	0.00
69	C	Pentachlorophenol	40.000	40.435	-1.1	114	0.00
70		Phenanthrene	40.000	39.591	1.0	117	0.00
71		Anthracene	40.000	39.614	1.0	116	0.00
72		Carbazole	40.000	39.051	2.4	115	0.00
73		Di-n-butylphthalate	40.000	39.218	2.0	115	0.00
74	C	Fluoranthene	40.000	37.780	5.5	111	0.00
75	I	Chrysene-d12	20.000	20.000	0.0	102	0.00
76		Benzidine	40.000	42.318	-5.8	119	0.00
77		Pyrene	40.000	42.310	-5.8	111	0.00
78	S	Terphenyl-d14	80.000	87.695	-9.6	116	0.00
79		Butylbenzylphthalate	40.000	40.267	-0.7	102	0.00
80		Benzo(a)anthracene	40.000	39.212	2.0	100	0.00
81		3,3'-Dichlorobenzidine	40.000	39.217	2.0	97	0.00
82		Chrysene	40.000	38.833	2.9	99	0.00
83		Bis(2-ethylhexyl)phthalate	40.000	39.001	2.5	98	0.00
84	c	Di-n-octyl phthalate	40.000	37.951	5.1	92	0.00
85		Indeno(1,2,3-cd)pyrene	40.000	37.755	5.6	85	0.00
86	I	Perylene-d12	20.000	20.000	0.0	90	0.00
87		Benzo(b)fluoranthene	40.000	38.863	2.8	87	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN072618\
Data File : BN002167.D
Acq On : 26 Jul 2018 10:08
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Jul 27 00:56:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN072518.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 26 18:27:25 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(k)fluoranthene	40.000	41.204	-3.0	94	0.00
89 C	Benzo(a)pyrene	40.000	40.050	-0.1	89	0.00
90	Dibenzo(a,h)anthracene	40.000	40.213	-0.5	85	0.00
91	Benzo(g,h,i)perylene	40.000	40.408	-1.0	85	0.00

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