Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN032719\
Data File : BN005002.D
Acq On : 27 Mar 2019 19:51
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Mar 28 04:50:51 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	104	0.00
2		1,4-Dioxane	40.000	38.484	3.8	101	0.00
3		Pyridine	40.000	38.792	3.0	102	0.00
4		n-Nitrosodimethylamine	40.000	39.718	0.7	101	0.00
5	S	2-Fluorophenol	80.000	77.002	3.7	101	0.00
6		Aniline	40.000	38.481	3.8	98	0.00
7	S	Phenol-d6	80.000	77.269	3.4	99	0.00
8		2-Chlorophenol	40.000	38.421	3.9	99	0.00
9		Benzaldehyde	40.000	40.260	-0.6	100	0.00
10	C	Phenol	40.000	38.321	4.2	98	0.00
11		bis(2-Chloroethyl)ether	40.000	38.688	3.3	99	0.00
12		1,3-Dichlorobenzene	40.000	38.379	4.1	99	0.00
13	C	1,4-Dichlorobenzene	40.000	38.830	2.9	101	0.00
14		1,2-Dichlorobenzene	40.000	38.613	3.5	99	0.00
15		Benzyl Alcohol	40.000	38.310	4.2	98	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.785	3.0	100	-0.01
17		2-Methylphenol	40.000	38.845	2.9	98	0.00
18		Hexachloroethane	40.000	38.303	4.2	99	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	39.445	1.4	99	0.00
20		3+4-Methylphenols	40.000	38.690	3.3	99	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	101	0.00
22		Acetophenone	40.000	39.600	1.0	99	0.00
23	S	Nitrobenzene-d5	80.000	78.150	2.3	98	0.00
24		Nitrobenzene	40.000	39.162	2.1	99	0.00
25		Isophorone	40.000	39.238	1.9	98	0.00
26	C	2-Nitrophenol	40.000	39.027	2.4	98	0.00
27		2,4-Dimethylphenol	40.000	38.920	2.7	98	0.00
28		bis(2-Chloroethoxy)methane	40.000	39.602	1.0	99	0.00
29	C	2,4-Dichlorophenol	40.000	39.365	1.6	99	0.00
30		1,2,4-Trichlorobenzene	40.000	39.797	0.5	101	0.00
31		Naphthalene	40.000	39.380	1.5	99	0.00
32		Benzoic acid	40.000	35.159	12.1	85	0.00
33		4-Chloroaniline	40.000	39.879	0.3	99	0.00
34	C	Hexachlorobutadiene	40.000	39.450	1.4	100	0.00
35		Caprolactam	40.000	38.997	2.5	93	0.00
36	C	4-Chloro-3-methylphenol	40.000	39.311	1.7	97	0.00
37		2-Methylnaphthalene	40.000	39.549	1.1	99	0.00
38		1-Methylnaphthalene	40.000	39.639	0.9	99	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	99	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.101	-0.3	101	0.00
41	P	Hexachlorocyclopentadiene	40.000	39.824	0.4	101	0.00
42	S	2,4,6-Tribromophenol	80.000	78.746	1.6	96	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.741	0.6	98	0.00
44		2,4,5-Trichlorophenol	40.000	39.302	1.7	96	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN032719\
Data File : BN005002.D
Acq On : 27 Mar 2019 19:51
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Mar 28 04:50:51 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	80.613	-0.8	100	0.00
46		1,1'-Biphenyl	40.000	40.452	-1.1	100	0.00
47		2-Chloronaphthalene	40.000	40.127	-0.3	100	0.00
48		2-Nitroaniline	40.000	39.336	1.7	96	0.00
49		Acenaphthylene	40.000	39.855	0.4	98	0.00
50		Dimethylphthalate	40.000	39.878	0.3	98	0.00
51		2,6-Dinitrotoluene	40.000	39.663	0.8	98	0.00
52	C	Acenaphthene	40.000	39.691	0.8	98	0.00
53		3-Nitroaniline	40.000	39.214	2.0	95	0.00
54	P	2,4-Dinitrophenol	40.000	37.485	6.3	90	0.00
55		Dibenzofuran	40.000	39.864	0.3	99	0.00
56	P	4-Nitrophenol	40.000	38.191	4.5	93	0.00
57		2,4-Dinitrotoluene	40.000	39.681	0.8	96	0.00
58		Fluorene	40.000	40.082	-0.2	98	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.761	0.6	97	0.00
60		Diethylphthalate	40.000	39.781	0.5	97	0.00
61		4-Chlorophenyl-phenylether	40.000	40.673	-1.7	100	0.00
62		4-Nitroaniline	40.000	38.910	2.7	93	0.00
63		Azobenzene	40.000	39.515	1.2	97	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	96	0.00
65		4,6-Dinitro-2-methylphenol	40.000	38.377	4.1	92	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.944	0.1	98	0.00
67		4-Bromophenyl-phenylether	40.000	39.814	0.5	97	0.00
68		Hexachlorobenzene	40.000	40.366	-0.9	98	0.00
69		Atrazine	40.000	38.913	2.7	92	0.00
70	C	Pentachlorophenol	40.000	38.440	3.9	95	0.00
71		Phenanthrene	40.000	39.294	1.8	95	0.00
72		Anthracene	40.000	39.517	1.2	95	0.00
73		Carbazole	40.000	39.011	2.5	93	0.00
74		Di-n-butylphthalate	40.000	38.576	3.6	92	0.00
75	C	Fluoranthene	40.000	38.026	4.9	91	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	91	0.00
77		Benzidine	40.000	39.295	1.8	80	0.00
78		Pyrene	40.000	41.322	-3.3	91	0.00
79	S	Terphenyl-d14	80.000	84.094	-5.1	91	0.00
80		Butylbenzylphthalate	40.000	38.130	4.7	84	0.00
81		Benzo(a)anthracene	40.000	38.951	2.6	87	0.00
82		3,3'-Dichlorobenzidine	40.000	39.256	1.9	87	0.00
83		Chrysene	40.000	39.096	2.3	89	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	38.052	4.9	85	0.00
85	c	Di-n-octyl phthalate	40.000	36.872	7.8	84	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	38.800	3.0	96	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	92	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA N\Data\BN032719\
Data File : BN005002.D
Acq On : 27 Mar 2019 19:51
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Mar 28 04:50:51 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	39.757	0.6	92	0.00
89	Benzo(k)fluoranthene	40.000	39.524	1.2	89	0.00
90 C	Benzo(a)pyrene	40.000	39.248	1.9	91	-0.01
91	Dibenzo(a,h)anthracene	40.000	40.189	-0.5	96	-0.01
92	Benzo(g,h,i)perylene	40.000	39.946	0.1	96	-0.01

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