Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052819\
Data File : BM020573.D
Acq On : 28 May 2019 10:11
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 28 13:20:13 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM052419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 24 16:07:05 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	110	-0.01
2		1,4-Dioxane	40.000	41.095	-2.7	112	0.00
3		Pyridine	40.000	39.365	1.6	106	0.00
4		n-Nitrosodimethylamine	40.000	42.671	-6.7	121	0.00
5	S	2-Fluorophenol	80.000	77.920	2.6	105	-0.01
6		Aniline	40.000	37.437	6.4	101	-0.01
7	S	Phenol-d6	80.000	76.415	4.5	103	-0.01
8		2-Chlorophenol	40.000	39.120	2.2	107	-0.01
9		Benzaldehyde	40.000	39.125	2.2	104	-0.01
10	C	Phenol	40.000	38.716	3.2	105	-0.02
11		bis(2-Chloroethyl)ether	40.000	37.206	7.0	102	-0.01
12		1,3-Dichlorobenzene	40.000	39.427	1.4	107	-0.01
13	C	1,4-Dichlorobenzene	40.000	38.786	3.0	105	-0.01
14		1,2-Dichlorobenzene	40.000	38.986	2.5	107	-0.01
15		Benzyl Alcohol	40.000	35.137	12.2	95	-0.01
16		2,2'-oxybis(1-Chloropropane	40.000	38.448	3.9	107	-0.01
17		2-Methylphenol	40.000	35.828	10.4	99	-0.01
18		Hexachloroethane	40.000	40.479	-1.2	111	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	36.594	8.5	99	-0.01
20		3+4-Methylphenols	40.000	35.538	11.2	97	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	98	0.00
22		Acetophenone	40.000	39.493	1.3	98	-0.01
23	S	Nitrobenzene-d5	80.000	80.613	-0.8	100	-0.01
24		Nitrobenzene	40.000	39.827	0.4	98	-0.01
25		Isophorone	40.000	39.672	0.8	98	-0.01
26	C	2-Nitrophenol	40.000	37.821	5.4	93	0.00
27		2,4-Dimethylphenol	40.000	38.178	4.6	96	-0.01
28		bis(2-Chloroethoxy)methane	40.000	38.251	4.4	95	-0.01
29	C	2,4-Dichlorophenol	40.000	38.488	3.8	95	-0.01
30		1,2,4-Trichlorobenzene	40.000	40.331	-0.8	102	-0.01
31		Naphthalene	40.000	39.696	0.8	99	-0.01
32		Benzoic acid	40.000	33.407	16.5	78	-0.03
33		4-Chloroaniline	40.000	35.358	11.6	86	-0.01
34	C	Hexachlorobutadiene	40.000	42.804	-7.0	107	-0.01
35		Caprolactam	40.000	35.651	10.9	88	-0.01
36	C	4-Chloro-3-methylphenol	40.000	36.126	9.7	89	-0.02
37		2-Methylnaphthalene	40.000	37.497	6.3	94	0.00
38		1-Methylnaphthalene	40.000	37.806	5.5	94	-0.01
39	I	Acenaphthene-d10	20.000	20.000	0.0	93	-0.01
40		1,2,4,5-Tetrachlorobenzene	40.000	41.226	-3.1	97	-0.01
41	P	Hexachlorocyclopentadiene	40.000	43.222	-8.1	108	0.00
42	S	2,4,6-Tribromophenol	80.000	73.294	8.4	87	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.049	2.4	90	-0.01
44		2,4,5-Trichlorophenol	40.000	36.993	7.5	85	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052819\
Data File : BM020573.D
Acq On : 28 May 2019 10:11
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 28 13:20:13 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM052419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 24 16:07:05 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	80.155	-0.2	95	0.00
46		1,1'-Biphenyl	40.000	39.439	1.4	93	-0.01
47		2-Chloronaphthalene	40.000	39.752	0.6	94	0.00
48		2-Nitroaniline	40.000	35.879	10.3	85	-0.01
49		Acenaphthylene	40.000	38.717	3.2	92	-0.01
50		Dimethylphthalate	40.000	38.413	4.0	91	-0.01
51		2,6-Dinitrotoluene	40.000	38.013	5.0	89	0.00
52	C	Acenaphthene	40.000	38.528	3.7	91	-0.01
53		3-Nitroaniline	40.000	34.102	14.7	79	0.00
54	P	2,4-Dinitrophenol	40.000	35.827	10.4	86	0.00
55		Dibenzofuran	40.000	38.265	4.3	91	-0.01
56	P	4-Nitrophenol	40.000	34.838	12.9	82	0.00
57		2,4-Dinitrotoluene	40.000	36.306	9.2	85	0.00
58		Fluorene	40.000	37.649	5.9	89	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	37.147	7.1	87	-0.01
60		Diethylphthalate	40.000	38.291	4.3	91	0.00
61		4-Chlorophenyl-phenylether	40.000	38.430	3.9	92	0.00
62		4-Nitroaniline	40.000	32.473	18.8	79	-0.01
63		Azobenzene	40.000	38.873	2.8	91	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	86	0.00
65		4,6-Dinitro-2-methylphenol	40.000	36.409	9.0	79	-0.01
66	c	n-Nitrosodiphenylamine	40.000	40.627	-1.6	89	0.00
67		4-Bromophenyl-phenylether	40.000	40.911	-2.3	88	-0.01
68		Hexachlorobenzene	40.000	41.051	-2.6	90	0.00
69		Atrazine	40.000	41.466	-3.7	87	-0.01
70	C	Pentachlorophenol	40.000	38.936	2.7	83	-0.01
71		Phenanthrene	40.000	39.198	2.0	86	0.00
72		Anthracene	40.000	38.050	4.9	83	0.00
73		Carbazole	40.000	36.309	9.2	79	0.00
74		Di-n-butylphthalate	40.000	39.807	0.5	87	-0.01
75	C	Fluoranthene	40.000	36.636	8.4	81	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	73	-0.01
77		Benzidine	40.000	34.887	12.8	57	0.00
78		Pyrene	40.000	44.649	-11.6	80	0.00
79	S	Terphenyl-d14	80.000	90.568	-13.2	81	0.00
80		Butylbenzylphthalate	40.000	43.295	-8.2	78	0.00
81		Benzo(a)anthracene	40.000	39.218	2.0	72	0.00
82		3,3'-Dichlorobenzidine	40.000	37.638	5.9	69	0.00
83		Chrysene	40.000	39.572	1.1	73	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	42.479	-6.2	79	0.00
85	c	Di-n-octyl phthalate	40.000	39.877	0.3	75	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	37.830	5.4	79	-0.02
87	I	Perylene-d12	20.000	20.000	0.0	71	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM052819\
Data File : BM020573.D
Acq On : 28 May 2019 10:11
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 28 13:20:13 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM052419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 24 16:07:05 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	38.155	4.6	69	-0.01
89	Benzo(k)fluoranthene	40.000	40.120	-0.3	69	0.00
90 C	Benzo(a)pyrene	40.000	39.184	2.0	71	0.00
91	Dibenzo(a,h)anthracene	40.000	40.812	-2.0	79	-0.02
92	Benzo(g,h,i)perylene	40.000	41.819	-4.5	82	-0.02

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