Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030519\
Data File : BM019052.D
Acq On : 06 Mar 2019 01:50
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 06 06:40:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 05 03:36:13 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	85	0.00
2		1,4-Dioxane	40.000	44.523	-11.3	95	0.00
3		Pyridine	40.000	46.488	-16.2	90	0.00
4		n-Nitrosodimethylamine	40.000	49.188	-23.0	100	0.00
5	S	2-Fluorophenol	80.000	90.191	-12.7	94	0.00
6		Aniline	40.000	43.097	-7.7	89	0.00
7	S	Phenol-d6	80.000	87.176	-9.0	90	0.00
8		2-Chlorophenol	40.000	45.142	-12.9	94	0.00
9		Benzaldehyde	40.000	49.222	-23.1	99	0.00
10	C	Phenol	40.000	43.906	-9.8	91	0.00
11		bis(2-Chloroethyl)ether	40.000	43.472	-8.7	90	0.00
12		1,3-Dichlorobenzene	40.000	44.495	-11.2	94	-0.01
13	C	1,4-Dichlorobenzene	40.000	44.455	-11.1	94	0.00
14		1,2-Dichlorobenzene	40.000	44.115	-10.3	93	0.00
15		Benzyl Alcohol	40.000	42.937	-7.3	87	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	44.158	-10.4	95	-0.01
17		2-Methylphenol	40.000	43.103	-7.8	89	0.00
18		Hexachloroethane	40.000	44.985	-12.5	94	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	42.558	-6.4	87	0.00
20		3+4-Methylphenols	40.000	43.061	-7.7	88	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	83	0.00
22		Acetophenone	40.000	42.102	-5.3	87	0.00
23	S	Nitrobenzene-d5	80.000	85.786	-7.2	87	0.00
24		Nitrobenzene	40.000	42.439	-6.1	87	0.00
25		Isophorone	40.000	44.687	-11.7	90	0.00
26	C	2-Nitrophenol	40.000	45.281	-13.2	90	0.00
27		2,4-Dimethylphenol	40.000	44.057	-10.1	90	0.00
28		bis(2-Chloroethoxy)methane	40.000	43.690	-9.2	89	0.00
29	C	2,4-Dichlorophenol	40.000	44.423	-11.1	89	0.00
30		1,2,4-Trichlorobenzene	40.000	43.409	-8.5	90	0.00
31		Naphthalene	40.000	43.810	-9.5	91	0.00
32		Benzoic acid	40.000	43.429	-8.6	89	-0.02
33		4-Chloroaniline	40.000	42.098	-5.2	85	0.00
34	C	Hexachlorobutadiene	40.000	42.805	-7.0	90	0.00
35		Caprolactam	40.000	39.099	2.3	78	-0.01
36	C	4-Chloro-3-methylphenol	40.000	42.659	-6.6	86	0.00
37		2-Methylnaphthalene	40.000	43.856	-9.6	91	0.00
38		1-Methylnaphthalene	40.000	43.650	-9.1	91	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	81	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	42.590	-6.5	87	0.00
41	P	Hexachlorocyclopentadiene	40.000	36.275	9.3	74	0.00
42	S	2,4,6-Tribromophenol	80.000	89.900	-12.4	89	0.00
43	C	2,4,6-Trichlorophenol	40.000	42.783	-7.0	84	0.00
44		2,4,5-Trichlorophenol	40.000	42.562	-6.4	84	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030519\
Data File : BM019052.D
Acq On : 06 Mar 2019 01:50
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 06 06:40:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 05 03:36:13 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	85.820	-7.3	88	-0.01
46		1,1'-Biphenyl	40.000	42.679	-6.7	88	0.00
47		2-Chloronaphthalene	40.000	42.194	-5.5	86	0.00
48		2-Nitroaniline	40.000	43.897	-9.7	85	0.00
49		Acenaphthylene	40.000	44.158	-10.4	89	0.00
50		Dimethylphthalate	40.000	41.966	-4.9	86	0.00
51		2,6-Dinitrotoluene	40.000	43.588	-9.0	86	0.00
52	C	Acenaphthene	40.000	42.403	-6.0	87	0.00
53		3-Nitroaniline	40.000	40.505	-1.3	82	0.00
54	P	2,4-Dinitrophenol	40.000	38.949	2.6	83	0.00
55		Dibenzofuran	40.000	42.144	-5.4	86	-0.01
56	P	4-Nitrophenol	40.000	40.727	-1.8	82	0.00
57		2,4-Dinitrotoluene	40.000	44.708	-11.8	87	0.00
58		Fluorene	40.000	43.972	-9.9	90	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.817	-4.5	83	0.00
60		Diethylphthalate	40.000	42.605	-6.5	87	0.00
61		4-Chlorophenyl-phenylether	40.000	41.290	-3.2	85	0.00
62		4-Nitroaniline	40.000	42.426	-6.1	85	0.00
63		Azobenzene	40.000	42.925	-7.3	86	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	83	0.00
65		4,6-Dinitro-2-methylphenol	40.000	39.075	2.3	81	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.766	-4.4	86	0.00
67		4-Bromophenyl-phenylether	40.000	41.581	-4.0	86	0.00
68		Hexachlorobenzene	40.000	41.923	-4.8	88	-0.01
69		Atrazine	40.000	41.123	-2.8	85	0.00
70	C	Pentachlorophenol	40.000	40.823	-2.1	80	0.00
71		Phenanthrene	40.000	42.542	-6.4	89	0.00
72		Anthracene	40.000	43.467	-8.7	90	0.00
73		Carbazole	40.000	42.589	-6.5	87	0.00
74		Di-n-butylphthalate	40.000	45.132	-12.8	91	0.00
75	C	Fluoranthene	40.000	43.192	-8.0	89	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	82	0.00
77		Benzidine	40.000	52.744	-31.9#	88	0.00
78		Pyrene	40.000	43.959	-9.9	89	0.00
79	S	Terphenyl-d14	80.000	87.218	-9.0	86	0.00
80		Butylbenzylphthalate	40.000	46.958	-17.4	91	0.00
81		Benzo(a)anthracene	40.000	42.886	-7.2	86	0.00
82		3,3'-Dichlorobenzidine	40.000	43.838	-9.6	86	0.00
83		Chrysene	40.000	42.843	-7.1	87	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	47.727	-19.3	94	0.00
85	c	Di-n-octyl phthalate	40.000	48.407	-21.0#	94	-0.01
86		Indeno(1,2,3-cd)pyrene	40.000	40.622	-1.6	76	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	72	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030519\
Data File : BM019052.D
Acq On : 06 Mar 2019 01:50
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 06 06:40:24 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 05 03:36:13 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	45.539	-13.8	83	0.00
89	Benzo(k)fluoranthene	40.000	42.390	-6.0	78	-0.01
90 C	Benzo(a)pyrene	40.000	42.928	-7.3	77	-0.01
91	Dibenzo(a,h)anthracene	40.000	44.356	-10.9	76	-0.01
92	Benzo(g,h,i)perylene	40.000	44.426	-11.1	75	-0.01

(#) = Out of Range SPCC's out = 0 CCC's out = 1