Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM042219\
Data File : BM020005.D
Acq On : 22 Apr 2019 14:39
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 22 17:27:05 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM041519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 19 00:56:08 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	107	-0.01
2		1,4-Dioxane	40.000	45.118	-12.8	122	0.00
3		Pyridine	40.000	43.533	-8.8	123	-0.01
4		n-Nitrosodimethylamine	40.000	40.973	-2.4	107	0.00
5	S	2-Fluorophenol	80.000	83.292	-4.1	111	-0.01
6		Aniline	40.000	35.940	10.2	96	0.00
7	S	Phenol-d6	80.000	74.575	6.8	100	-0.01
8		2-Chlorophenol	40.000	38.659	3.4	104	-0.01
9		Benzaldehyde	40.000	37.611	6.0	100	-0.01
10	C	Phenol	40.000	37.076	7.3	100	-0.01
11		bis(2-Chloroethyl)ether	40.000	37.009	7.5	98	-0.01
12		1,3-Dichlorobenzene	40.000	41.234	-3.1	111	0.00
13	C	1,4-Dichlorobenzene	40.000	40.943	-2.4	110	-0.01
14		1,2-Dichlorobenzene	40.000	39.455	1.4	106	-0.01
15		Benzyl Alcohol	40.000	34.891	12.8	93	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	34.723	13.2	92	-0.01
17		2-Methylphenol	40.000	35.011	12.5	93	-0.01
18		Hexachloroethane	40.000	40.892	-2.2	109	-0.01
19	P	n-Nitroso-di-n-propylamine	40.000	32.118	19.7	85	0.00
20		3+4-Methylphenols	40.000	33.963	15.1	90	-0.01
21	I	Naphthalene-d8	20.000	20.000	0.0	89	-0.01
22		Acetophenone	40.000	40.067	-0.2	90	-0.01
23	S	Nitrobenzene-d5	80.000	83.577	-4.5	93	-0.01
24		Nitrobenzene	40.000	41.586	-4.0	92	-0.01
25		Isophorone	40.000	36.581	8.5	81	-0.01
26	C	2-Nitrophenol	40.000	39.545	1.1	88	0.00
27		2,4-Dimethylphenol	40.000	38.381	4.0	85	0.00
28		bis(2-Chloroethoxy)methane	40.000	38.040	4.9	85	0.00
29	C	2,4-Dichlorophenol	40.000	39.146	2.1	87	-0.01
30		1,2,4-Trichlorobenzene	40.000	42.471	-6.2	95	-0.01
31		Naphthalene	40.000	40.155	-0.4	89	-0.01
32		Benzoic acid	40.000	20.723	48.2#	41	-0.02
33		4-Chloroaniline	40.000	37.012	7.5	82	-0.01
34	C	Hexachlorobutadiene	40.000	44.576	-11.4	100	-0.02
35		Caprolactam	40.000	33.256	16.9	74	0.00
36	C	4-Chloro-3-methylphenol	40.000	35.438	11.4	79	0.00
37		2-Methylnaphthalene	40.000	37.386	6.5	84	-0.01
38		1-Methylnaphthalene	40.000	37.066	7.3	83	-0.01
39	I	Acenaphthene-d10	20.000	20.000	0.0	78	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	43.691	-9.2	86	-0.01
41	P	Hexachlorocyclopentadiene	40.000	60.799	-52.0#	150	-0.01
42	S	2,4,6-Tribromophenol	80.000	75.402	5.7	75	-0.01
43	C	2,4,6-Trichlorophenol	40.000	40.228	-0.6	79	-0.01
44		2,4,5-Trichlorophenol	40.000	38.481	3.8	75	-0.01

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM042219\
Data File : BM020005.D
Acq On : 22 Apr 2019 14:39
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 22 17:27:05 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM041519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 19 00:56:08 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	83.538	-4.4	82	-0.01
46		1,1'-Biphenyl	40.000	41.523	-3.8	82	0.00
47		2-Chloronaphthalene	40.000	41.465	-3.7	81	-0.01
48		2-Nitroaniline	40.000	38.679	3.3	75	0.00
49		Acenaphthylene	40.000	39.202	2.0	77	0.00
50		Dimethylphthalate	40.000	38.744	3.1	76	0.00
51		2,6-Dinitrotoluene	40.000	38.577	3.6	76	0.00
52	C	Acenaphthene	40.000	39.883	0.3	79	-0.01
53		3-Nitroaniline	40.000	36.770	8.1	72	0.00
54	P	2,4-Dinitrophenol	40.000	36.457	8.9	72	0.00
55		Dibenzofuran	40.000	38.901	2.7	77	-0.01
56	P	4-Nitrophenol	40.000	37.646	5.9	75	-0.02
57		2,4-Dinitrotoluene	40.000	36.808	8.0	73	0.00
58		Fluorene	40.000	37.557	6.1	75	-0.01
59		2,3,4,6-Tetrachlorophenol	40.000	37.883	5.3	75	0.00
60		Diethylphthalate	40.000	38.140	4.6	75	0.00
61		4-Chlorophenyl-phenylether	40.000	38.974	2.6	77	0.00
62		4-Nitroaniline	40.000	35.350	11.6	68	0.00
63		Azobenzene	40.000	39.569	1.1	77	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	74	-0.01
65		4,6-Dinitro-2-methylphenol	40.000	33.507	16.2	62	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.452	1.4	74	-0.01
67		4-Bromophenyl-phenylether	40.000	40.087	-0.2	76	0.00
68		Hexachlorobenzene	40.000	40.321	-0.8	76	-0.01
69		Atrazine	40.000	32.621	18.4	59	-0.01
70	C	Pentachlorophenol	40.000	35.285	11.8	66	0.00
71		Phenanthrene	40.000	38.874	2.8	73	0.00
72		Anthracene	40.000	38.775	3.1	73	0.00
73		Carbazole	40.000	38.290	4.3	72	0.00
74		Di-n-butylphthalate	40.000	39.638	0.9	74	0.00
75	C	Fluoranthene	40.000	38.778	3.1	74	-0.01
76	I	Chrysene-d12	20.000	20.000	0.0	73	0.00
77		Benzidine	40.000	42.111	-5.3	77	-0.01
78		Pyrene	40.000	40.969	-2.4	74	0.00
79	S	Terphenyl-d14	80.000	80.266	-0.3	73	0.00
80		Butylbenzylphthalate	40.000	41.577	-3.9	75	-0.01
81		Benzo(a)anthracene	40.000	42.882	-7.2	79	0.00
82		3,3'-Dichlorobenzidine	40.000	44.514	-11.3	80	0.00
83		Chrysene	40.000	42.808	-7.0	79	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	41.859	-4.6	77	0.00
85	c	Di-n-octyl phthalate	40.000	44.808	-12.0	83	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	63.695	-59.2#	118	-0.01
87	I	Perylene-d12	20.000	20.000	0.0	96	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM042219\
Data File : BM020005.D
Acq On : 22 Apr 2019 14:39
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 22 17:27:05 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM041519.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 19 00:56:08 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	37.879	5.3	93	0.00
89	Benzo(k)fluoranthene	40.000	38.609	3.5	93	0.00
90 C	Benzo(a)pyrene	40.000	40.117	-0.3	98	0.00
91	Dibenzo(a,h)anthracene	40.000	48.309	-20.8	117	-0.01
92	Benzo(g,h,i)perylene	40.000	49.870	-24.7	120	-0.02

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