Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM031919\
Data File : BM019229.D
Acq On : 19 Mar 2019 10:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 19 12:20:03 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 19 12:12:53 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	81	-0.02
2		1,4-Dioxane	40.000	34.427	13.9	73	-0.02
3		Pyridine	40.000	39.394	1.5	76	-0.02
4		n-Nitrosodimethylamine	40.000	34.617	13.5	70	-0.02
5	S	2-Fluorophenol	80.000	75.167	6.0	77	-0.02
6		Aniline	40.000	39.414	1.5	80	-0.02
7	S	Phenol-d6	80.000	78.957	1.3	79	-0.02
8		2-Chlorophenol	40.000	38.674	3.3	79	-0.03
9		Benzaldehyde	40.000	36.728	8.2	74	-0.02
10	C	Phenol	40.000	39.703	0.7	80	-0.02
11		bis(2-Chloroethyl)ether	40.000	38.527	3.7	78	-0.02
12		1,3-Dichlorobenzene	40.000	37.758	5.6	78	-0.03
13	C	1,4-Dichlorobenzene	40.000	37.485	6.3	78	-0.02
14		1,2-Dichlorobenzene	40.000	38.130	4.7	79	-0.02
15		Benzyl Alcohol	40.000	39.689	0.8	79	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	37.918	5.2	78	-0.03
17		2-Methylphenol	40.000	39.788	0.5	80	-0.02
18		Hexachloroethane	40.000	37.612	6.0	76	-0.02
19	P	n-Nitroso-di-n-propylamine	40.000	40.976	-2.4	82	-0.02
20		3+4-Methylphenols	40.000	41.714	-4.3	83	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	85	-0.03
22		Acetophenone	40.000	38.171	4.6	81	-0.03
23	S	Nitrobenzene-d5	80.000	75.861	5.2	81	-0.02
24		Nitrobenzene	40.000	37.859	5.4	81	-0.02
25		Isophorone	40.000	39.157	2.1	83	-0.02
26	C	2-Nitrophenol	40.000	40.108	-0.3	82	-0.03
27		2,4-Dimethylphenol	40.000	39.163	2.1	83	-0.02
28		bis(2-Chloroethoxy)methane	40.000	38.706	3.2	83	-0.02
29	C	2,4-Dichlorophenol	40.000	40.166	-0.4	83	-0.02
30		1,2,4-Trichlorobenzene	40.000	37.680	5.8	82	-0.03
31		Naphthalene	40.000	37.528	6.2	82	-0.02
32		Benzoic acid	40.000	37.338	6.7	79	-0.03
33		4-Chloroaniline	40.000	39.022	2.4	82	-0.02
34	C	Hexachlorobutadiene	40.000	37.350	6.6	81	-0.03
35		Caprolactam	40.000	42.837	-7.1	88	-0.02
36	C	4-Chloro-3-methylphenol	40.000	39.778	0.6	84	-0.03
37		2-Methylnaphthalene	40.000	38.344	4.1	84	-0.03
38		1-Methylnaphthalene	40.000	38.153	4.6	83	-0.02
39	I	Acenaphthene-d10	20.000	20.000	0.0	88	-0.02
40		1,2,4,5-Tetrachlorobenzene	40.000	37.520	6.2	84	-0.02
41	P	Hexachlorocyclopentadiene	40.000	28.382	29.0#	62	-0.02
42	S	2,4,6-Tribromophenol	80.000	78.258	2.2	85	-0.02
43	C	2,4,6-Trichlorophenol	40.000	39.517	1.2	86	-0.02
44		2,4,5-Trichlorophenol	40.000	39.321	1.7	85	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM031919\
Data File : BM019229.D
Acq On : 19 Mar 2019 10:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 19 12:20:03 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 19 12:12:53 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	74.408	7.0	84	-0.02
46		1,1'-Biphenyl	40.000	37.162	7.1	84	-0.02
47		2-Chloronaphthalene	40.000	37.637	5.9	85	-0.02
48		2-Nitroaniline	40.000	39.189	2.0	82	-0.02
49		Acenaphthylene	40.000	37.869	5.3	84	-0.02
50		Dimethylphthalate	40.000	37.749	5.6	85	-0.02
51		2,6-Dinitrotoluene	40.000	39.022	2.4	85	-0.02
52	C	Acenaphthene	40.000	37.379	6.6	84	-0.02
53		3-Nitroaniline	40.000	37.159	7.1	83	-0.02
54	P	2,4-Dinitrophenol	40.000	31.957	20.1	70	-0.02
55		Dibenzofuran	40.000	37.449	6.4	85	-0.02
56	P	4-Nitrophenol	40.000	36.208	9.5	80	-0.02
57		2,4-Dinitrotoluene	40.000	37.654	5.9	85	-0.02
58		Fluorene	40.000	37.617	6.0	84	-0.02
59		2,3,4,6-Tetrachlorophenol	40.000	38.415	4.0	84	-0.02
60		Diethylphthalate	40.000	38.044	4.9	85	-0.02
61		4-Chlorophenyl-phenylether	40.000	37.872	5.3	85	-0.02
62		4-Nitroaniline	40.000	36.936	7.7	83	-0.02
63		Azobenzene	40.000	38.128	4.7	84	-0.02
64	I	Phenanthrene-d10	20.000	20.000	0.0	90	-0.02
65		4,6-Dinitro-2-methylphenol	40.000	34.676	13.3	77	-0.02
66	c	n-Nitrosodiphenylamine	40.000	38.242	4.4	86	-0.02
67		4-Bromophenyl-phenylether	40.000	38.568	3.6	87	-0.02
68		Hexachlorobenzene	40.000	38.298	4.3	87	-0.02
69		Atrazine	40.000	38.314	4.2	85	-0.02
70	C	Pentachlorophenol	40.000	36.076	9.8	81	-0.02
71		Phenanthrene	40.000	37.138	7.2	85	-0.02
72		Anthracene	40.000	37.953	5.1	85	-0.02
73		Carbazole	40.000	37.150	7.1	83	-0.02
74		Di-n-butylphthalate	40.000	39.724	0.7	87	-0.02
75	C	Fluoranthene	40.000	36.778	8.1	84	-0.02
76	I	Chrysene-d12	20.000	20.000	0.0	83	-0.02
77		Benzidine	40.000	31.916	20.2	65	-0.02
78		Pyrene	40.000	40.798	-2.0	82	-0.02
79	S	Terphenyl-d14	80.000	83.197	-4.0	84	-0.02
80		Butylbenzylphthalate	40.000	43.746	-9.4	85	-0.02
81		Benzo(a)anthracene	40.000	37.208	7.0	78	-0.02
82		3,3'-Dichlorobenzidine	40.000	37.810	5.5	77	-0.02
83		Chrysene	40.000	37.354	6.6	78	-0.02
84		Bis(2-ethylhexyl)phthalate	40.000	43.880	-9.7	87	-0.02
85	c	Di-n-octyl phthalate	40.000	41.142	-2.9	85	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	33.180	17.1	78	-0.04
87	I	Perylene-d12	20.000	20.000	0.0	80	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM031919\
Data File : BM019229.D
Acq On : 19 Mar 2019 10:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 19 12:20:03 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\8270-BM031119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Mar 19 12:12:53 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	37.494	6.3	76	-0.02
89	Benzo(k)fluoranthene	40.000	39.060	2.3	76	-0.02
90 C	Benzo(a)pyrene	40.000	37.582	6.0	76	-0.03
91	Dibenzo(a,h)anthracene	40.000	37.163	7.1	78	-0.04
92	Benzo(g,h,i)perylene	40.000	38.038	4.9	80	-0.04

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