Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM042319\
Data File : BM020053.D
Acq On : 24 Apr 2019 02:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 24 02:59:39 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	87	-0.02
2		1,4-Dioxane	40.000	42.519	-6.3	93	-0.02
3		Pyridine	40.000	39.389	1.5	90	-0.02
4		n-Nitrosodimethylamine	40.000	38.628	3.4	82	-0.01
5	S	2-Fluorophenol	80.000	75.229	6.0	82	-0.02
6		Aniline	40.000	34.347	14.1	75	-0.02
7	S	Phenol-d6	80.000	70.787	11.5	77	-0.02
8		2-Chlorophenol	40.000	36.611	8.5	80	-0.02
9		Benzaldehyde	40.000	47.473	-18.7	96	-0.02
10	C	Phenol	40.000	35.025	12.4	77	-0.02
11		bis(2-Chloroethyl)ether	40.000	34.502	13.7	74	-0.02
12		1,3-Dichlorobenzene	40.000	38.495	3.8	84	-0.02
13	C	1,4-Dichlorobenzene	40.000	39.139	2.2	86	-0.02
14		1,2-Dichlorobenzene	40.000	37.903	5.2	83	-0.02
15		Benzyl Alcohol	40.000	34.175	14.6	74	-0.02
16		2,2'-oxybis(1-Chloropropane	40.000	31.633	20.9	69	-0.02
17		2-Methylphenol	40.000	33.811	15.5	73	-0.02
18		Hexachloroethane	40.000	39.461	1.3	86	-0.03
19	P	n-Nitroso-di-n-propylamine	40.000	31.317	21.7	68	-0.02
20		3+4-Methylphenols	40.000	32.280	19.3	70	-0.02
21	I	Naphthalene-d8	20.000	20.000	0.0	76	-0.02
22		Acetophenone	40.000	37.315	6.7	72	-0.02
23	S	Nitrobenzene-d5	80.000	78.544	1.8	75	-0.02
24		Nitrobenzene	40.000	39.032	2.4	74	-0.02
25		Isophorone	40.000	34.261	14.3	65	-0.02
26	C	2-Nitrophenol	40.000	35.950	10.1	68	-0.02
27		2,4-Dimethylphenol	40.000	35.494	11.3	68	-0.02
28		bis(2-Chloroethoxy)methane	40.000	35.303	11.7	68	-0.02
29	C	2,4-Dichlorophenol	40.000	37.147	7.1	70	-0.02
30		1,2,4-Trichlorobenzene	40.000	40.021	-0.1	76	-0.02
31		Naphthalene	40.000	37.910	5.2	72	-0.02
32		Benzoic acid	40.000	27.415	31.5#	50	0.00
33		4-Chloroaniline	40.000	34.647	13.4	66	-0.02
34	C	Hexachlorobutadiene	40.000	42.934	-7.3	83	-0.03
35		Caprolactam	40.000	31.081	22.3	59	-0.02
36	C	4-Chloro-3-methylphenol	40.000	34.072	14.8	65	-0.02
37		2-Methylnaphthalene	40.000	36.075	9.8	69	-0.02
38		1-Methylnaphthalene	40.000	35.969	10.1	69	-0.02
39	I	Acenaphthene-d10	20.000	20.000	0.0	70	-0.02
40		1,2,4,5-Tetrachlorobenzene	40.000	40.907	-2.3	73	-0.02
41	P	Hexachlorocyclopentadiene	40.000	43.845	-9.6	89	-0.03
42	S	2,4,6-Tribromophenol	80.000	71.039	11.2	64	-0.02
43	C	2,4,6-Trichlorophenol	40.000	37.436	6.4	67	-0.02
44		2,4,5-Trichlorophenol	40.000	35.183	12.0	62	-0.02

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM042319\
Data File : BM020053.D
Acq On : 24 Apr 2019 02:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 24 02:59:39 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	77.516	3.1	69	-0.03
46		1,1'-Biphenyl	40.000	38.142	4.6	68	-0.02
47		2-Chloronaphthalene	40.000	38.228	4.4	68	-0.02
48		2-Nitroaniline	40.000	36.175	9.6	63	-0.02
49		Acenaphthylene	40.000	36.848	7.9	66	-0.02
50		Dimethylphthalate	40.000	36.603	8.5	65	-0.02
51		2,6-Dinitrotoluene	40.000	35.402	11.5	63	-0.02
52	C	Acenaphthene	40.000	37.228	6.9	67	-0.02
53		3-Nitroaniline	40.000	34.161	14.6	60	-0.02
54	P	2,4-Dinitrophenol	40.000	35.130	12.2	62	-0.01
55		Dibenzofuran	40.000	36.947	7.6	66	-0.02
56	P	4-Nitrophenol	40.000	29.936	25.2#	51	-0.02
57		2,4-Dinitrotoluene	40.000	34.650	13.4	62	-0.02
58		Fluorene	40.000	36.468	8.8	65	-0.02
59		2,3,4,6-Tetrachlorophenol	40.000	36.655	8.4	66	-0.02
60		Diethylphthalate	40.000	36.172	9.6	64	-0.02
61		4-Chlorophenyl-phenylether	40.000	37.439	6.4	67	-0.02
62		4-Nitroaniline	40.000	33.167	17.1	58	-0.02
63		Azobenzene	40.000	38.033	4.9	67	-0.02
64	I	Phenanthrene-d10	20.000	20.000	0.0	69	-0.02
65		4,6-Dinitro-2-methylphenol	40.000	31.102	22.2	54	-0.02
66	c	n-Nitrosodiphenylamine	40.000	37.012	7.5	65	-0.02
67		4-Bromophenyl-phenylether	40.000	37.518	6.2	66	-0.02
68		Hexachlorobenzene	40.000	37.620	6.0	66	-0.02
69		Atrazine	40.000	34.438	13.9	58	-0.02
70	C	Pentachlorophenol	40.000	41.739	-4.3	73	-0.02
71		Phenanthrene	40.000	36.877	7.8	65	-0.02
72		Anthracene	40.000	36.614	8.5	65	-0.02
73		Carbazole	40.000	36.287	9.3	64	-0.02
74		Di-n-butylphthalate	40.000	35.922	10.2	63	-0.02
75	C	Fluoranthene	40.000	37.268	6.8	66	-0.02
76	I	Chrysene-d12	20.000	20.000	0.0	73	-0.02
77		Benzidine	40.000	42.468	-6.2	77	-0.02
78		Pyrene	40.000	37.079	7.3	67	-0.02
79	S	Terphenyl-d14	80.000	73.151	8.6	67	-0.02
80		Butylbenzylphthalate	40.000	35.645	10.9	64	-0.02
81		Benzo(a)anthracene	40.000	39.859	0.4	74	-0.02
82		3,3'-Dichlorobenzidine	40.000	40.784	-2.0	73	-0.02
83		Chrysene	40.000	39.802	0.5	73	-0.02
84		Bis(2-ethylhexyl)phthalate	40.000	35.420	11.4	65	-0.02
85	c	Di-n-octyl phthalate	40.000	38.603	3.5	71	-0.02
86		Indeno(1,2,3-cd)pyrene	40.000	58.701	-46.8#	108	-0.04
87	I	Perylene-d12	20.000	20.000	0.0	94	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM042319\
Data File : BM020053.D
Acq On : 24 Apr 2019 02:25
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Apr 24 02:59:39 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	35.706	10.7	86	-0.02
89	Benzo(k)fluoranthene	40.000	35.487	11.3	84	-0.02
90 C	Benzo(a)pyrene	40.000	37.236	6.9	89	-0.03
91	Dibenzo(a,h)anthracene	40.000	45.245	-13.1	108	-0.05
92	Benzo(g,h,i)perylene	40.000	46.380	-16.0	110	-0.05

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