Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM053019\
Data File : BM020612.D
Acq On : 29 May 2019 23:28
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 30 02:45:20 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM052419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 30 02:15:06 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	73	0.00
2		1,4-Dioxane	0.533	0.711	-33.4#	96	0.00
3		Pyridine	1.511	1.801	-19.2	82	0.00
4		n-Nitrosodimethylamine	0.378	0.394	-4.2	78	0.00
5	S	2-Fluorophenol	1.019	1.108	-8.7	78	0.00
6		Aniline	2.226	2.154	3.2	69	0.00
7	S	Phenol-d6	1.539	1.605	-4.3	75	0.00
8		2-Chlorophenol	1.247	1.325	-6.3	77	0.00
9		Benzaldehyde	0.982	0.994	-1.2	71	0.00
10	C	Phenol	1.662	1.724	-3.7	74	0.00
11		bis(2-Chloroethyl)ether	1.240	1.196	3.5	70	0.00
12		1,3-Dichlorobenzene	1.596	1.729	-8.3	78	0.00
13	C	1,4-Dichlorobenzene	1.596	1.709	-7.1	77	0.00
14		1,2-Dichlorobenzene	1.600	1.742	-8.9	79	0.00
15		Benzyl Alcohol	1.424	1.321	7.2	67	0.00
16		2,2'-oxybis(1-Chloropropane	0.864	0.825	4.5	70	0.00
17		2-Methylphenol	1.097	1.054	3.9	70	0.00
18		Hexachloroethane	0.646	0.741	-14.7	83	0.00
19	P	n-Nitroso-di-n-propylamine	1.366	1.304	4.5	68	0.00
20		3+4-Methylphenols	1.471	1.398	5.0	68	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	65	0.00
22		Acetophenone	0.551	0.566	-2.7	67	0.00
23	S	Nitrobenzene-d5	0.476	0.535	-12.4	74	0.00
24		Nitrobenzene	0.471	0.517	-9.8	72	0.00
25		Isophorone	0.887	0.944	-6.4	70	0.00
26	C	2-Nitrophenol	0.174	0.169	2.9	63	0.00
27		2,4-Dimethylphenol	0.259	0.256	1.2	66	0.00
28		bis(2-Chloroethoxy)methane	0.462	0.452	2.2	64	0.00
29	C	2,4-Dichlorophenol	0.344	0.370	-7.6	71	0.00
30		1,2,4-Trichlorobenzene	0.470	0.544	-15.7	77	0.00
31		Naphthalene	1.033	1.124	-8.8	72	0.00
32		Benzoic acid	0.154	0.078	49.4#	32#	0.02
33		4-Chloroaniline	0.412	0.419	-1.7	66	0.00
34	C	Hexachlorobutadiene	0.350	0.455	-30.0#	86	0.00
35		Caprolactam	0.113	0.104	8.0	60	0.00
36	C	4-Chloro-3-methylphenol	0.374	0.360	3.7	63	0.00
37		2-Methylnaphthalene	0.789	0.831	-5.3	70	0.00
38		1-Methylnaphthalene	0.764	0.801	-4.8	69	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	65	0.00
40		1,2,4,5-Tetrachlorobenzene	0.841	0.978	-16.3	76	0.00
41	P	Hexachlorocyclopentadiene	0.342	0.152	55.6#	29#	0.00
42	S	2,4,6-Tribromophenol	0.401	0.436	-8.7	72	0.00
43	C	2,4,6-Trichlorophenol	0.520	0.565	-8.7	70	0.00
44		2,4,5-Trichlorophenol	0.486	0.515	-6.0	68	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM053019\
Data File : BM020612.D
Acq On : 29 May 2019 23:28
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 30 02:45:20 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM052419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 30 02:15:06 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.594	1.756	-10.2	73	0.00
46		1,1'-Biphenyl	1.540	1.645	-6.8	70	0.00
47		2-Chloronaphthalene	1.306	1.434	-9.8	72	0.00
48		2-Nitroaniline	0.384	0.355	7.6	60	0.00
49		Acenaphthylene	1.986	2.059	-3.7	69	0.00
50		Dimethylphthalate	1.772	1.857	-4.8	69	0.00
51		2,6-Dinitrotoluene	0.369	0.379	-2.7	67	0.00
52	C	Acenaphthene	1.197	1.270	-6.1	70	0.00
53		3-Nitroaniline	0.309	0.260	15.9	54	0.00
54	P	2,4-Dinitrophenol	0.187	0.143	23.5	49#	0.00
55		Dibenzofuran	2.071	2.233	-7.8	71	0.00
56	P	4-Nitrophenol	0.199	0.201	-1.0	64	0.00
57		2,4-Dinitrotoluene	0.533	0.515	3.4	63	0.00
58		Fluorene	1.720	1.785	-3.8	68	0.00
59		2,3,4,6-Tetrachlorophenol	0.565	0.605	-7.1	70	0.00
60		Diethylphthalate	1.745	1.795	-2.9	68	0.00
61		4-Chlorophenyl-phenylether	1.098	1.187	-8.1	72	0.00
62		4-Nitroaniline	0.310	0.248	20.0	54	0.00
63		Azobenzene	1.698	1.828	-7.7	70	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	62	0.00
65		4,6-Dinitro-2-methylphenol	0.118	0.080	32.2#	41#	0.00
66	c	n-Nitrosodiphenylamine	0.542	0.582	-7.4	67	0.00
67		4-Bromophenyl-phenylether	0.266	0.298	-12.0	69	0.00
68		Hexachlorobenzene	0.323	0.384	-18.9	75	0.00
69		Atrazine	0.223	0.217	2.7	58	0.00
70	C	Pentachlorophenol	0.167	0.182	-9.0	66	0.00
71		Phenanthrene	1.087	1.160	-6.7	67	0.00
72		Anthracene	1.054	1.069	-1.4	63	0.00
73		Carbazole	0.904	0.876	3.1	60	0.00
74		Di-n-butylphthalate	1.112	1.151	-3.5	65	0.00
75	C	Fluoranthene	1.511	1.568	-3.8	66	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	61	0.00
77		Benzidine	0.368	0.298	19.0	44#	0.00
78		Pyrene	1.256	1.393	-10.9	67	0.00
79	S	Terphenyl-d14	1.106	1.225	-10.8	67	0.00
80		Butylbenzylphthalate	0.441	0.419	5.0	57	0.00
81		Benzo(a)anthracene	1.384	1.459	-5.4	65	0.00
82		3,3'-Dichlorobenzidine	0.505	0.476	5.7	58	0.00
83		Chrysene	1.338	1.433	-7.1	66	0.00
84		Bis(2-ethylhexyl)phthalate	0.643	0.632	1.7	61	0.00
85	c	Di-n-octyl phthalate	1.118	1.091	2.4	61	0.00
86		Indeno(1,2,3-cd)pyrene	1.617	1.802	-11.4	78	-0.02
87	I	Perylene-d12	1.000	1.000	0.0	63	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM053019\
Data File : BM020612.D
Acq On : 29 May 2019 23:28
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 30 02:45:20 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM052419.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu May 30 02:15:06 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.309	1.342	-2.5	66	0.00
89	Benzo(k)fluoranthene	1.270	1.423	-12.0	69	0.00
90 C	Benzo(a)pyrene	1.205	1.290	-7.1	68	0.00
91	Dibenzo(a,h)anthracene	1.233	1.373	-11.4	76	-0.01
92	Benzo(g,h,i)perylene	1.143	1.313	-14.9	80	0.00

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