Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF081519\
Data File : BF116271.D
Acq On : 15 Aug 2019 11:32
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 15 15:05:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 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	105	0.00
2		1,4-Dioxane	40.000	41.050	-2.6	110	-0.02
3		Pyridine	40.000	37.749	5.6	100	-0.02
4		n-Nitrosodimethylamine	40.000	37.193	7.0	98	-0.02
5	S	2-Fluorophenol	80.000	87.240	-9.0	113	0.00
6		Aniline	40.000	39.675	0.8	102	0.00
7	S	Phenol-d6	80.000	84.107	-5.1	110	0.00
8		2-Chlorophenol	40.000	43.991	-10.0	114	0.00
9		Benzaldehyde	40.000	38.224	4.4	99	0.00
10	C	Phenol	40.000	41.486	-3.7	108	0.00
11		bis(2-Chloroethyl)ether	40.000	43.124	-7.8	112	0.00
12		1,3-Dichlorobenzene	40.000	41.713	-4.3	108	0.00
13	C	1,4-Dichlorobenzene	40.000	42.390	-6.0	110	0.00
14		1,2-Dichlorobenzene	40.000	42.023	-5.1	107	0.00
15		Benzyl Alcohol	40.000	39.384	1.5	100	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.192	-3.0	106	0.00
17		2-Methylphenol	40.000	42.308	-5.8	112	0.00
18		Hexachloroethane	40.000	42.315	-5.8	109	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	40.844	-2.1	108	0.00
20		3+4-Methylphenols	40.000	42.546	-6.4	108	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	107	0.00
22		Acetophenone	40.000	40.745	-1.9	107	0.00
23	S	Nitrobenzene-d5	80.000	81.499	-1.9	108	0.00
24		Nitrobenzene	40.000	42.219	-5.5	112	0.00
25		Isophorone	40.000	43.130	-7.8	115	0.00
26	C	2-Nitrophenol	40.000	44.042	-10.1	116	0.00
27		2,4-Dimethylphenol	40.000	41.706	-4.3	110	0.00
28		bis(2-Chloroethoxy)methane	40.000	43.011	-7.5	114	0.00
29	C	2,4-Dichlorophenol	40.000	44.014	-10.0	114	0.00
30		1,2,4-Trichlorobenzene	40.000	41.872	-4.7	110	0.00
31		Naphthalene	40.000	42.070	-5.2	109	0.00
32		Benzoic acid	40.000	38.874	2.8	113	0.01
33		4-Chloroaniline	40.000	40.407	-1.0	106	0.00
34	C	Hexachlorobutadiene	40.000	41.506	-3.8	109	0.00
35		Caprolactam	40.000	43.495	-8.7	114	0.01
36	C	4-Chloro-3-methylphenol	40.000	43.969	-9.9	116	0.00
37		2-Methylnaphthalene	40.000	41.806	-4.5	109	0.00
38		1-Methylnaphthalene	40.000	41.961	-4.9	110	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	109	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	41.522	-3.8	111	0.00
41	P	Hexachlorocyclopentadiene	40.000	36.139	9.7	98	0.00
42	S	2,4,6-Tribromophenol	80.000	84.468	-5.6	113	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.205	2.0	105	0.00
44		2,4,5-Trichlorophenol	40.000	39.756	0.6	107	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF081519\
Data File : BF116271.D
Acq On : 15 Aug 2019 11:32
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 15 15:05:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 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	77.774	2.8	103	0.00
46		1,1'-Biphenyl	40.000	40.846	-2.1	109	0.00
47		2-Chloronaphthalene	40.000	39.967	0.1	107	0.00
48		2-Nitroaniline	40.000	40.686	-1.7	110	0.00
49		Acenaphthylene	40.000	40.594	-1.5	108	0.00
50		Dimethylphthalate	40.000	41.318	-3.3	111	0.00
51		2,6-Dinitrotoluene	40.000	41.787	-4.5	112	0.00
52	C	Acenaphthene	40.000	40.492	-1.2	108	0.00
53		3-Nitroaniline	40.000	40.869	-2.2	111	0.00
54	P	2,4-Dinitrophenol	40.000	37.475	6.3	104	0.00
55		Dibenzofuran	40.000	38.728	3.2	102	0.00
56	P	4-Nitrophenol	40.000	38.931	2.7	104	0.00
57		2,4-Dinitrotoluene	40.000	38.481	3.8	102	0.00
58		Fluorene	40.000	41.729	-4.3	111	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	39.041	2.4	105	0.00
60		Diethylphthalate	40.000	41.866	-4.7	111	0.00
61		4-Chlorophenyl-phenylether	40.000	41.725	-4.3	110	0.00
62		4-Nitroaniline	40.000	38.540	3.7	103	0.00
63		Azobenzene	40.000	41.676	-4.2	111	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	108	0.00
65		4,6-Dinitro-2-methylphenol	40.000	43.983	-10.0	115	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.881	-4.7	112	0.00
67		4-Bromophenyl-phenylether	40.000	41.806	-4.5	111	0.00
68		Hexachlorobenzene	40.000	41.015	-2.5	109	0.00
69		Atrazine	40.000	33.701	15.7	90	0.00
70	C	Pentachlorophenol	40.000	39.994	0.0	104	0.00
71		Phenanthrene	40.000	41.179	-2.9	109	0.00
72		Anthracene	40.000	41.426	-3.6	111	0.00
73		Carbazole	40.000	42.903	-7.3	114	0.00
74		Di-n-butylphthalate	40.000	43.697	-9.2	113	0.00
75	C	Fluoranthene	40.000	42.007	-5.0	112	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	107	0.00
77		Benzidine	40.000	39.367	1.6	98	0.00
78		Pyrene	40.000	41.026	-2.6	112	0.00
79	S	Terphenyl-d14	80.000	78.549	1.8	107	0.00
80		Butylbenzylphthalate	40.000	43.970	-9.9	117	0.00
81		Benzo(a)anthracene	40.000	42.917	-7.3	115	0.00
82		3,3'-Dichlorobenzidine	40.000	44.290	-10.7	116	0.00
83		Chrysene	40.000	41.898	-4.7	112	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	45.224	-13.1	119	0.00
85	c	Di-n-octyl phthalate	40.000	45.838	-14.6	119	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	48.571	-21.4	136	0.01
87	I	Perylene-d12	20.000	20.000	0.0	122	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA F\Data\BF081519\
Data File : BF116271.D
Acq On : 15 Aug 2019 11:32
Operator : HP/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 15 15:05:59 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF073019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Aug 07 11:13:18 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	38.699	3.3	112	0.00
89	Benzo(k)fluoranthene	40.000	40.336	-0.8	127	0.00
90 C	Benzo(a)pyrene	40.000	41.255	-3.1	125	0.00
91	Dibenzo(a,h)anthracene	40.000	43.678	-9.2	138	0.01
92	Benzo(g,h,i)perylene	40.000	43.999	-10.0	143	0.00

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