Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022620\
Data File : BF119087.D
Acq On : 26 Feb 2020 11:17
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 26 16:14:32 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 26 16:03:05 2020
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	132	0.00
2		1,4-Dioxane	40.000	35.815	10.5	125	0.00
3		Pyridine	40.000	37.352	6.6	131	0.00
4		n-Nitrosodimethylamine	40.000	40.824	-2.1	142	0.00
5	S	2-Fluorophenol	80.000	76.281	4.6	130	0.00
6		Aniline	40.000	39.024	2.4	132	0.00
7	S	Phenol-d6	80.000	78.802	1.5	134	0.00
8		2-Chlorophenol	40.000	39.731	0.7	135	0.00
9		Benzaldehyde	40.000	34.617	13.5	119	0.00
10	C	Phenol	40.000	38.682	3.3	132	0.00
11		bis(2-Chloroethyl)ether	40.000	40.592	-1.5	140	0.00
12		1,3-Dichlorobenzene	40.000	39.300	1.8	136	0.00
13	C	1,4-Dichlorobenzene	40.000	39.332	1.7	135	0.00
14		1,2-Dichlorobenzene	40.000	39.136	2.2	133	0.00
15		Benzyl Alcohol	40.000	41.366	-3.4	139	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.667	-4.2	145	0.00
17		2-Methylphenol	40.000	40.754	-1.9	140	0.00
18		Hexachloroethane	40.000	41.154	-2.9	143	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	43.005	-7.5	150	0.00
20		3+4-Methylphenols	40.000	38.943	2.6	137	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	138	0.00
22		Acetophenone	40.000	39.797	0.5	145	0.00
23	S	Nitrobenzene-d5	80.000	78.497	1.9	143	0.00
24		Nitrobenzene	40.000	38.961	2.6	143	0.00
25		Isophorone	40.000	41.142	-2.9	152	0.00
26	C	2-Nitrophenol	40.000	39.954	0.1	146	0.00
27		2,4-Dimethylphenol	40.000	39.240	1.9	143	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.025	-0.1	146	0.00
29	C	2,4-Dichlorophenol	40.000	39.768	0.6	143	0.00
30		1,2,4-Trichlorobenzene	40.000	38.706	3.2	140	0.00
31		Naphthalene	40.000	38.151	4.6	138	0.00
32		Benzoic acid	40.000	44.406	-11.0	174	0.00
33		4-Chloroaniline	40.000	37.995	5.0	137	0.00
34	C	Hexachlorobutadiene	40.000	39.479	1.3	144	0.00
35		Caprolactam	40.000	41.925	-4.8	150	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.844	-2.1	148	0.00
37		2-Methylnaphthalene	40.000	39.081	2.3	142	0.00
38		1-Methylnaphthalene	40.000	39.149	2.1	141	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	143	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	37.649	5.9	144	0.00
41	P	Hexachlorocyclopentadiene	40.000	38.637	3.4	157	0.00
42	S	2,4,6-Tribromophenol	80.000	82.845	-3.6	158	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.272	1.8	150	0.00
44		2,4,5-Trichlorophenol	40.000	38.701	3.2	146	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022620\
Data File : BF119087.D
Acq On : 26 Feb 2020 11:17
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 26 16:14:32 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 26 16:03:05 2020
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	72.042	9.9	141	0.00
46		1,1'-Biphenyl	40.000	37.636	5.9	142	0.00
47		2-Chloronaphthalene	40.000	38.292	4.3	146	0.00
48		2-Nitroaniline	40.000	40.292	-0.7	153	0.00
49		Acenaphthylene	40.000	37.599	6.0	142	0.00
50		Dimethylphthalate	40.000	40.835	-2.1	157	0.00
51		2,6-Dinitrotoluene	40.000	40.261	-0.7	153	0.00
52	C	Acenaphthene	40.000	38.087	4.8	143	0.00
53		3-Nitroaniline	40.000	39.266	1.8	148	0.00
54	P	2,4-Dinitrophenol	40.000	41.967	-4.9	168	0.00
55		Dibenzofuran	40.000	36.885	7.8	137	0.00
56	P	4-Nitrophenol	40.000	35.949	10.1	133	0.00
57		2,4-Dinitrotoluene	40.000	38.067	4.8	141	0.00
58		Fluorene	40.000	38.385	4.0	144	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	40.348	-0.9	155	0.00
60		Diethylphthalate	40.000	41.278	-3.2	156	0.00
61		4-Chlorophenyl-phenylether	40.000	38.964	2.6	146	0.00
62		4-Nitroaniline	40.000	39.655	0.9	149	0.00
63		Azobenzene	40.000	39.951	0.1	152	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	145	0.00
65		4,6-Dinitro-2-methylphenol	40.000	42.431	-6.1	164	0.00
66	c	n-Nitrosodiphenylamine	40.000	39.356	1.6	151	0.00
67		4-Bromophenyl-phenylether	40.000	39.437	1.4	153	0.00
68		Hexachlorobenzene	40.000	39.544	1.1	153	0.00
69		Atrazine	40.000	37.748	5.6	146	0.00
70	C	Pentachlorophenol	40.000	40.118	-0.3	157	0.00
71		Phenanthrene	40.000	37.479	6.3	144	0.00
72		Anthracene	40.000	37.711	5.7	142	0.00
73		Carbazole	40.000	37.963	5.1	144	0.00
74		Di-n-butylphthalate	40.000	41.239	-3.1	158	0.00
75	C	Fluoranthene	40.000	38.041	4.9	144	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	132	0.00
77		Benzidine	40.000	33.608	16.0	115	0.00
78		Pyrene	40.000	38.825	2.9	141	0.00
79	S	Terphenyl-d14	80.000	78.485	1.9	144	0.00
80		Butylbenzylphthalate	40.000	43.923	-9.8	156	0.00
81		Benzo(a)anthracene	40.000	38.856	2.9	136	0.00
82		3,3'-Dichlorobenzidine	40.000	39.740	0.6	134	0.00
83		Chrysene	40.000	38.553	3.6	136	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	45.443	-13.6	160	0.00
85	c	Di-n-octyl phthalate	40.000	45.158	-12.9	151	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	35.171	12.1	96	0.00
87	I	Perylene-d12	20.000	20.000	0.0	86	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA F\DATA\BF022620\
Data File : BF119087.D
Acq On : 26 Feb 2020 11:17
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 26 16:14:32 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA F\METHODS\8270-BF022020.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 26 16:03:05 2020
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	39.751	0.6	129	0.00
89	Benzo(k)fluoranthene	40.000	41.564	-3.9	123	0.00
90 C	Benzo(a)pyrene	40.000	38.649	3.4	90	0.00
91	Dibenzo(a,h)anthracene	40.000	38.652	3.4	95	0.00
92	Benzo(g,h,i)perylene	40.000	37.759	5.6	94	0.00

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