Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051819\
Data File : BF114404.D
Acq On : 18 May 2019 21:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 20 07:39:12 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 10 15:56:46 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	100	0.00
2		1,4-Dioxane	40.000	33.475	16.3	82	0.00
3		Pyridine	40.000	34.809	13.0	89	0.00
4		n-Nitrosodimethylamine	40.000	35.923	10.2	90	0.00
5	S	2-Fluorophenol	80.000	73.442	8.2	90	0.00
6		Aniline	40.000	36.400	9.0	90	0.00
7	S	Phenol-d6	80.000	77.113	3.6	97	0.00
8		2-Chlorophenol	40.000	39.872	0.3	99	0.00
9		Benzaldehyde	40.000	35.161	12.1	83	0.00
10	C	Phenol	40.000	36.678	8.3	90	0.00
11		bis(2-Chloroethyl)ether	40.000	40.935	-2.3	102	0.00
12		1,3-Dichlorobenzene	40.000	39.330	1.7	99	0.00
13	C	1,4-Dichlorobenzene	40.000	39.139	2.2	98	0.00
14		1,2-Dichlorobenzene	40.000	39.125	2.2	96	0.00
15		Benzyl Alcohol	40.000	37.266	6.8	92	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	38.002	5.0	94	0.00
17		2-Methylphenol	40.000	38.117	4.7	96	0.00
18		Hexachloroethane	40.000	39.437	1.4	98	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	38.257	4.4	98	0.00
20		3+4-Methylphenols	40.000	39.179	2.1	98	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	101	0.00
22		Acetophenone	40.000	40.211	-0.5	100	0.00
23	S	Nitrobenzene-d5	80.000	80.067	-0.1	99	0.00
24		Nitrobenzene	40.000	40.567	-1.4	100	0.00
25		Isophorone	40.000	40.863	-2.2	106	0.00
26	C	2-Nitrophenol	40.000	42.475	-6.2	107	0.00
27		2,4-Dimethylphenol	40.000	37.116	7.2	93	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.566	-1.4	103	0.00
29	C	2,4-Dichlorophenol	40.000	40.404	-1.0	100	0.00
30		1,2,4-Trichlorobenzene	40.000	39.855	0.4	99	0.00
31		Naphthalene	40.000	39.631	0.9	96	0.00
32		Benzoic acid	40.000	32.488	18.8	82	0.00
33		4-Chloroaniline	40.000	40.560	-1.4	99	0.00
34	C	Hexachlorobutadiene	40.000	39.346	1.6	97	0.00
35		Caprolactam	40.000	44.526	-11.3	107	0.00
36	C	4-Chloro-3-methylphenol	40.000	42.118	-5.3	102	0.00
37		2-Methylnaphthalene	40.000	40.801	-2.0	99	0.00
38		1-Methylnaphthalene	40.000	41.274	-3.2	99	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	105	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	39.881	0.3	100	0.00
41	P	Hexachlorocyclopentadiene	40.000	18.118	54.7#	40	0.00
42	S	2,4,6-Tribromophenol	80.000	70.715	11.6	94	0.00
43	C	2,4,6-Trichlorophenol	40.000	39.198	2.0	99	0.00
44		2,4,5-Trichlorophenol	40.000	38.635	3.4	97	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051819\
Data File : BF114404.D
Acq On : 18 May 2019 21:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 20 07:39:12 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 10 15:56:46 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	72.266	9.7	96	0.00
46		1,1'-Biphenyl	40.000	38.837	2.9	99	0.00
47		2-Chloronaphthalene	40.000	38.469	3.8	98	0.00
48		2-Nitroaniline	40.000	40.472	-1.2	104	0.00
49		Acenaphthylene	40.000	38.030	4.9	97	0.00
50		Dimethylphthalate	40.000	40.980	-2.4	106	0.00
51		2,6-Dinitrotoluene	40.000	39.907	0.2	103	0.00
52	C	Acenaphthene	40.000	37.575	6.1	97	0.00
53		3-Nitroaniline	40.000	41.174	-2.9	105	0.00
54	P	2,4-Dinitrophenol	40.000	25.031	37.4#	62	0.00
55		Dibenzofuran	40.000	36.080	9.8	94	0.00
56	P	4-Nitrophenol	40.000	29.398	26.5#	78	0.00
57		2,4-Dinitrotoluene	40.000	36.233	9.4	95	0.00
58		Fluorene	40.000	38.032	4.9	100	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	36.042	9.9	94	0.00
60		Diethylphthalate	40.000	38.932	2.7	104	0.00
61		4-Chlorophenyl-phenylether	40.000	38.279	4.3	100	0.00
62		4-Nitroaniline	40.000	39.028	2.4	105	0.00
63		Azobenzene	40.000	38.593	3.5	102	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	100	0.00
65		4,6-Dinitro-2-methylphenol	40.000	33.535	16.2	82	0.00
66	c	n-Nitrosodiphenylamine	40.000	41.356	-3.4	99	0.00
67		4-Bromophenyl-phenylether	40.000	40.898	-2.2	100	0.00
68		Hexachlorobenzene	40.000	39.447	1.4	97	0.00
69		Atrazine	40.000	38.770	3.1	96	0.00
70	C	Pentachlorophenol	40.000	38.455	3.9	91	0.00
71		Phenanthrene	40.000	39.584	1.0	96	0.00
72		Anthracene	40.000	39.595	1.0	95	0.00
73		Carbazole	40.000	39.066	2.3	94	0.00
74		Di-n-butylphthalate	40.000	43.949	-9.9	106	0.00
75	C	Fluoranthene	40.000	36.258	9.4	88	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	74	0.00
77		Benzidine	40.000	31.427	21.4	60	0.00
78		Pyrene	40.000	44.122	-10.3	85	0.00
79	S	Terphenyl-d14	80.000	91.162	-14.0	88	0.00
80		Butylbenzylphthalate	40.000	52.276	-30.7#	97	0.00
81		Benzo(a)anthracene	40.000	41.772	-4.4	76	0.00
82		3,3'-Dichlorobenzidine	40.000	42.679	-6.7	75	0.00
83		Chrysene	40.000	40.591	-1.5	74	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	54.145	-35.4#	100	0.00
85	c	Di-n-octyl phthalate	40.000	49.051	-22.6#	88	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	53.856	-34.6#	103	0.00
87	I	Perylene-d12	20.000	20.000	0.0	85	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA_F\DATA\BF051819\
Data File : BF114404.D
Acq On : 18 May 2019 21:54
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: May 20 07:39:12 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_F\METHODS\8270-BF051019.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri May 10 15:56:46 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	39.849	0.4	85	0.00
89	Benzo(k)fluoranthene	40.000	38.463	3.8	77	0.00
90 C	Benzo(a)pyrene	40.000	41.850	-4.6	88	0.00
91	Dibenzo(a,h)anthracene	40.000	47.688	-19.2	103	0.00
92	Benzo(g,h,i)perylene	40.000	46.970	-17.4	104	0.00

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