Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP112719\
Data File : BP001093.D
Acq On : 27 Nov 2019 21:37
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Nov 28 03:03:18 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 27 17:38:46 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	92	0.00
2		1,4-Dioxane	40.000	42.264	-5.7	99	0.00
3		Pyridine	40.000	40.299	-0.7	95	0.00
4		n-Nitrosodimethylamine	40.000	41.340	-3.4	100	0.00
5	S	2-Fluorophenol	80.000	83.609	-4.5	97	0.00
6		Aniline	40.000	41.426	-3.6	99	0.00
7	S	Phenol-d6	80.000	82.296	-2.9	98	0.00
8		2-Chlorophenol	40.000	41.780	-4.5	98	0.00
9		Benzaldehyde	40.000	42.831	-7.1	98	0.00
10	C	Phenol	40.000	41.421	-3.6	97	0.00
11		bis(2-Chloroethyl)ether	40.000	41.622	-4.1	100	0.00
12		1,3-Dichlorobenzene	40.000	41.843	-4.6	99	0.00
13	C	1,4-Dichlorobenzene	40.000	41.431	-3.6	97	0.00
14		1,2-Dichlorobenzene	40.000	41.506	-3.8	97	0.00
15		Benzyl Alcohol	40.000	42.090	-5.2	100	0.00
16		2,2'-oxybis(1-Chloropropane	40.000	41.433	-3.6	99	-0.01
17		2-Methylphenol	40.000	40.721	-1.8	98	0.00
18		Hexachloroethane	40.000	41.576	-3.9	98	0.00
19	P	n-Nitroso-di-n-propylamine	40.000	41.313	-3.3	101	0.00
20		3+4-Methylphenols	40.000	40.974	-2.4	98	0.00
21	I	Naphthalene-d8	20.000	20.000	0.0	95	0.00
22		Acetophenone	40.000	40.338	-0.8	100	0.00
23	S	Nitrobenzene-d5	80.000	80.660	-0.8	99	0.00
24		Nitrobenzene	40.000	40.617	-1.5	99	0.00
25		Isophorone	40.000	40.366	-0.9	101	0.00
26	C	2-Nitrophenol	40.000	39.624	0.9	95	0.00
27		2,4-Dimethylphenol	40.000	40.604	-1.5	100	0.00
28		bis(2-Chloroethoxy)methane	40.000	40.485	-1.2	100	0.00
29	C	2,4-Dichlorophenol	40.000	41.040	-2.6	100	0.00
30		1,2,4-Trichlorobenzene	40.000	40.714	-1.8	99	0.00
31		Naphthalene	40.000	40.994	-2.5	99	0.00
32		Benzoic acid	40.000	41.962	-4.9	100	0.00
33		4-Chloroaniline	40.000	40.168	-0.4	100	0.00
34	C	Hexachlorobutadiene	40.000	40.870	-2.2	99	0.00
35		Caprolactam	40.000	40.002	-0.0	102	0.00
36	C	4-Chloro-3-methylphenol	40.000	40.547	-1.4	101	0.00
37		2-Methylnaphthalene	40.000	40.567	-1.4	99	0.00
38		1-Methylnaphthalene	40.000	41.001	-2.5	101	0.00
39	I	Acenaphthene-d10	20.000	20.000	0.0	98	0.00
40		1,2,4,5-Tetrachlorobenzene	40.000	40.655	-1.6	101	0.00
41	P	Hexachlorocyclopentadiene	40.000	40.759	-1.9	97	0.00
42	S	2,4,6-Tribromophenol	80.000	83.139	-3.9	103	0.00
43	C	2,4,6-Trichlorophenol	40.000	40.357	-0.9	100	0.00
44		2,4,5-Trichlorophenol	40.000	40.674	-1.7	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP112719\
Data File : BP001093.D
Acq On : 27 Nov 2019 21:37
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Nov 28 03:03:18 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 27 17:38:46 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	81.231	-1.5	100	0.00
46		1,1'-Biphenyl	40.000	40.927	-2.3	101	0.00
47		2-Chloronaphthalene	40.000	40.704	-1.8	101	0.00
48		2-Nitroaniline	40.000	40.293	-0.7	103	0.00
49		Acenaphthylene	40.000	41.125	-2.8	102	0.00
50		Dimethylphthalate	40.000	40.749	-1.9	103	0.00
51		2,6-Dinitrotoluene	40.000	40.431	-1.1	102	0.00
52	C	Acenaphthene	40.000	40.505	-1.3	102	0.00
53		3-Nitroaniline	40.000	40.518	-1.3	103	0.00
54	P	2,4-Dinitrophenol	40.000	35.782	10.5	94	0.00
55		Dibenzofuran	40.000	41.031	-2.6	102	0.00
56	P	4-Nitrophenol	40.000	41.057	-2.6	104	0.00
57		2,4-Dinitrotoluene	40.000	41.659	-4.1	101	0.00
58		Fluorene	40.000	41.266	-3.2	102	0.00
59		2,3,4,6-Tetrachlorophenol	40.000	41.466	-3.7	103	0.00
60		Diethylphthalate	40.000	41.155	-2.9	104	0.00
61		4-Chlorophenyl-phenylether	40.000	41.022	-2.6	103	0.00
62		4-Nitroaniline	40.000	40.496	-1.2	103	0.00
63		Azobenzene	40.000	40.802	-2.0	103	0.00
64	I	Phenanthrene-d10	20.000	20.000	0.0	100	0.00
65		4,6-Dinitro-2-methylphenol	40.000	37.561	6.1	99	0.00
66	c	n-Nitrosodiphenylamine	40.000	40.234	-0.6	103	0.00
67		4-Bromophenyl-phenylether	40.000	40.009	-0.0	101	0.00
68		Hexachlorobenzene	40.000	39.992	0.0	102	0.00
69		Atrazine	40.000	41.479	-3.7	104	0.00
70	C	Pentachlorophenol	40.000	41.146	-2.9	102	0.00
71		Phenanthrene	40.000	40.754	-1.9	104	0.00
72		Anthracene	40.000	40.613	-1.5	102	0.00
73		Carbazole	40.000	40.656	-1.6	103	0.00
74		Di-n-butylphthalate	40.000	42.916	-7.3	106	0.00
75	C	Fluoranthene	40.000	41.986	-5.0	103	0.00
76	I	Chrysene-d12	20.000	20.000	0.0	99	0.00
77		Benzidine	40.000	44.636	-11.6	92	0.00
78		Pyrene	40.000	39.229	1.9	104	0.00
79	S	Terphenyl-d14	80.000	79.126	1.1	103	0.00
80		Butylbenzylphthalate	40.000	41.536	-3.8	105	0.00
81		Benzo(a)anthracene	40.000	40.416	-1.0	103	0.00
82		3,3'-Dichlorobenzidine	40.000	43.203	-8.0	105	0.00
83		Chrysene	40.000	41.372	-3.4	104	0.00
84		Bis(2-ethylhexyl)phthalate	40.000	43.291	-8.2	107	0.00
85	c	Di-n-octyl phthalate	40.000	43.403	-8.5	107	0.00
86		Indeno(1,2,3-cd)pyrene	40.000	41.068	-2.7	105	0.00
87	I	Perylene-d12	20.000	20.000	0.0	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP112719\
Data File : BP001093.D
Acq On : 27 Nov 2019 21:37
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Nov 28 03:03:18 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP112719.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Nov 27 17:38:46 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	41.758	-4.4	109	0.00
89	Benzo(k)fluoranthene	40.000	38.440	3.9	100	0.00
90 C	Benzo(a)pyrene	40.000	40.387	-1.0	105	0.00
91	Dibenzo(a,h)anthracene	40.000	40.697	-1.7	105	0.00
92	Benzo(g,h,i)perylene	40.000	40.259	-0.6	105	0.00

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