Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120219\
Data File : BP001121.D
Acq On : 02 Dec 2019 12:45
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Dec 02 13:59:31 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
1	I	1,4-Dichlorobenzene-d4	1.000	1.000	0.0	94	0.00
2		1,4-Dioxane	0.563	0.580	-3.0	99	0.00
3		Pyridine	1.562	1.675	-7.2	103	0.00
4		n-Nitrosodimethylamine	0.676	0.743	-9.9	109	0.00
5	S	2-Fluorophenol	1.205	1.290	-7.1	101	0.00
6		Aniline	2.129	2.237	-5.1	103	0.00
7	S	Phenol-d6	1.606	1.715	-6.8	104	0.00
8		2-Chlorophenol	1.247	1.322	-6.0	102	0.00
9		Benzaldehyde	1.044	1.076	-3.1	108	0.00
10	C	Phenol	1.827	1.928	-5.5	102	0.00
11		bis(2-Chloroethyl)ether	1.423	1.504	-5.7	104	0.00
12		1,3-Dichlorobenzene	1.478	1.574	-6.5	103	0.00
13	C	1,4-Dichlorobenzene	1.489	1.574	-5.7	102	0.00
14		1,2-Dichlorobenzene	1.402	1.474	-5.1	101	0.00
15		Benzyl Alcohol	1.318	1.420	-7.7	105	0.00
16		2,2'-oxybis(1-Chloropropane	2.287	2.404	-5.1	103	0.00
17		2-Methylphenol	1.144	1.180	-3.1	102	0.00
18		Hexachloroethane	0.616	0.656	-6.5	103	0.00
19	P	n-Nitroso-di-n-propylamine	1.159	1.184	-2.2	102	0.00
20		3+4-Methylphenols	1.442	1.484	-2.9	101	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	95	0.00
22		Acetophenone	0.588	0.618	-5.1	103	0.00
23	S	Nitrobenzene-d5	0.461	0.491	-6.5	104	0.00
24		Nitrobenzene	0.458	0.487	-6.3	103	0.00
25		Isophorone	0.827	0.852	-3.0	103	0.00
26	C	2-Nitrophenol	0.168	0.173	-3.0	98	0.00
27		2,4-Dimethylphenol	0.266	0.280	-5.3	103	0.00
28		bis(2-Chloroethoxy)methane	0.519	0.533	-2.7	101	0.00
29	C	2,4-Dichlorophenol	0.303	0.316	-4.3	101	0.00
30		1,2,4-Trichlorobenzene	0.354	0.375	-5.9	103	0.00
31		Naphthalene	1.021	1.075	-5.3	102	0.00
32		Benzoic acid	0.192	0.191	0.5	94	0.01
33		4-Chloroaniline	0.443	0.458	-3.4	102	0.00
34	C	Hexachlorobutadiene	0.222	0.240	-8.1	104	0.00
35		Caprolactam	0.104	0.100	3.8	97	0.01
36	C	4-Chloro-3-methylphenol	0.359	0.369	-2.8	102	0.00
37		2-Methylnaphthalene	0.697	0.720	-3.3	100	0.00
38		1-Methylnaphthalene	0.658	0.676	-2.7	101	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
40		1,2,4,5-Tetrachlorobenzene	0.630	0.690	-9.5	103	0.00
41	P	Hexachlorocyclopentadiene	0.291	0.315	-8.2	97	0.00
42	S	2,4,6-Tribromophenol	0.192	0.196	-2.1	96	0.00
43	C	2,4,6-Trichlorophenol	0.412	0.438	-6.3	100	0.00
44		2,4,5-Trichlorophenol	0.426	0.448	-5.2	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120219\
Data File : BP001121.D
Acq On : 02 Dec 2019 12:45
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Dec 02 13:59:31 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.367	1.465	-7.2	99	0.00
46		1,1'-Biphenyl	1.546	1.667	-7.8	101	0.00
47		2-Chloronaphthalene	1.235	1.320	-6.9	100	0.00
48		2-Nitroaniline	0.484	0.507	-4.8	101	0.00
49		Acenaphthylene	1.851	1.964	-6.1	99	0.00
50		Dimethylphthalate	1.496	1.526	-2.0	97	0.00
51		2,6-Dinitrotoluene	0.320	0.326	-1.9	97	0.00
52	C	Acenaphthene	1.113	1.176	-5.7	100	0.00
53		3-Nitroaniline	0.360	0.360	0.0	96	0.00
54	P	2,4-Dinitrophenol	0.112	0.108	3.6	93	0.00
55		Dibenzofuran	1.673	1.764	-5.4	99	0.00
56	P	4-Nitrophenol	0.255	0.237	7.1	89	0.00
57		2,4-Dinitrotoluene	0.401	0.415	-3.5	95	0.00
58		Fluorene	1.340	1.384	-3.3	97	0.00
59		2,3,4,6-Tetrachlorophenol	0.351	0.365	-4.0	98	0.00
60		Diethylphthalate	1.549	1.544	0.3	96	0.00
61		4-Chlorophenyl-phenylether	0.683	0.714	-4.5	99	0.00
62		4-Nitroaniline	0.417	0.430	-3.1	99	0.00
63		Azobenzene	1.674	1.706	-1.9	97	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	86	0.00
65		4,6-Dinitro-2-methylphenol	0.080	0.071	11.3	79	0.00
66	c	n-Nitrosodiphenylamine	0.608	0.666	-9.5	97	0.00
67		4-Bromophenyl-phenylether	0.217	0.237	-9.2	95	0.00
68		Hexachlorobenzene	0.239	0.260	-8.8	96	0.00
69		Atrazine	0.186	0.186	0.0	87	0.00
70	C	Pentachlorophenol	0.124	0.123	0.8	85	0.00
71		Phenanthrene	1.051	1.119	-6.5	93	0.00
72		Anthracene	1.036	1.112	-7.3	93	0.00
73		Carbazole	0.943	0.981	-4.0	91	0.00
74		Di-n-butylphthalate	1.268	1.322	-4.3	88	0.00
75	C	Fluoranthene	1.113	1.167	-4.9	89	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	83	0.00
77		Benzidine	0.516	0.644	-24.8	86	0.00
78		Pyrene	1.575	1.583	-0.5	89	0.00
79	S	Terphenyl-d14	1.081	1.099	-1.7	89	0.00
80		Butylbenzylphthalate	0.728	0.753	-3.4	88	0.00
81		Benzo(a)anthracene	1.387	1.427	-2.9	89	0.00
82		3,3'-Dichlorobenzidine	0.458	0.487	-6.3	87	0.00
83		Chrysene	1.242	1.313	-5.7	90	0.00
84		Bis(2-ethylhexyl)phthalate	1.000	1.058	-5.8	88	0.00
85	c	Di-n-octyl phthalate	1.777	1.904	-7.1	89	0.00
86		Indeno(1,2,3-cd)pyrene	1.463	1.549	-5.9	91	0.00
87	I	Perylene-d12	1.000	1.000	0.0	85	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP120219\
Data File : BP001121.D
Acq On : 02 Dec 2019 12:45
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Dec 02 13:59:31 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	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.222	1.307	-7.0	93	0.00
89	Benzo(k)fluoranthene	1.177	1.193	-1.4	88	0.00
90 C	Benzo(a)pyrene	1.125	1.179	-4.8	91	0.00
91	Dibenzo(a,h)anthracene	1.101	1.148	-4.3	90	0.00
92	Benzo(g,h,i)perylene	1.087	1.128	-3.8	91	0.00

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