Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102418\
Data File : BN003288.D
Acq On : 25 Oct 2018 00:59
Operator : MJ/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Oct 25 05:26:26 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN102418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 24 14:21:25 2018
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	111	0.00
2		1,4-Dioxane	0.545	0.513	5.9	107	0.00
3		Pyridine	1.246	1.275	-2.3	107	0.00
4		n-Nitrosodimethylamine	0.459	0.443	3.5	108	0.00
5	S	2-Fluorophenol	1.148	1.122	2.3	107	0.00
6		Aniline	2.039	2.065	-1.3	110	0.00
7	S	Phenol-d6	1.633	1.636	-0.2	109	0.00
8		2-Chlorophenol	1.414	1.396	1.3	109	0.00
9		Benzaldehyde	0.948	0.920	3.0	107	0.00
10	C	Phenol	1.723	1.719	0.2	108	0.00
11		bis(2-Chloroethyl)ether	1.407	1.384	1.6	108	0.00
12		1,3-Dichlorobenzene	1.572	1.543	1.8	108	0.00
13	C	1,4-Dichlorobenzene	1.619	1.591	1.7	109	0.00
14		1,2-Dichlorobenzene	1.571	1.525	2.9	107	0.00
15		Benzyl Alcohol	1.177	1.188	-0.9	113	0.00
16		2,2'-oxybis(1-Chloropropane	2.063	2.029	1.6	110	0.00
17		2-Methylphenol	1.178	1.178	0.0	109	0.00
18		Hexachloroethane	0.567	0.553	2.5	109	0.00
19	P	n-Nitroso-di-n-propylamine	1.177	1.196	-1.6	111	0.00
20		3+4-Methylphenols	1.673	1.682	-0.5	109	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	113	0.00
22		Acetophenone	0.485	0.484	0.2	111	0.00
23	S	Nitrobenzene-d5	0.349	0.345	1.1	110	0.00
24		Nitrobenzene	0.356	0.352	1.1	110	0.00
25		Isophorone	0.612	0.616	-0.7	112	0.00
26	C	2-Nitrophenol	0.187	0.189	-1.1	110	0.00
27		2,4-Dimethylphenol	0.263	0.259	1.5	109	0.00
28		bis(2-Chloroethoxy)methane	0.412	0.411	0.2	112	0.00
29	C	2,4-Dichlorophenol	0.301	0.305	-1.3	111	0.00
30		1,2,4-Trichlorobenzene	0.342	0.334	2.3	110	0.00
31		Naphthalene	0.977	0.951	2.7	110	0.00
32		Benzoic acid	0.209	0.197	5.7	104	0.00
33		4-Chloroaniline	0.432	0.429	0.7	111	0.00
34	C	Hexachlorobutadiene	0.204	0.200	2.0	110	0.00
35		Caprolactam	0.118	0.121	-2.5	109	0.00
36	C	4-Chloro-3-methylphenol	0.335	0.339	-1.2	111	0.00
37		2-Methylnaphthalene	0.698	0.689	1.3	111	0.00
38		1-Methylnaphthalene	0.682	0.670	1.8	111	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
40		1,2,4,5-Tetrachlorobenzene	0.662	0.650	1.8	111	0.00
41	P	Hexachlorocyclopentadiene	0.229	0.242	-5.7	117	0.00
42	S	2,4,6-Tribromophenol	0.269	0.270	-0.4	113	0.00
43	C	2,4,6-Trichlorophenol	0.418	0.418	0.0	112	0.00
44		2,4,5-Trichlorophenol	0.436	0.438	-0.5	115	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102418\
Data File : BN003288.D
Acq On : 25 Oct 2018 00:59
Operator : MJ/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Oct 25 05:26:26 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN102418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 24 14:21:25 2018
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	1.474	1.426	3.3	111	0.00
46		1,1'-Biphenyl	1.755	1.705	2.8	112	0.00
47		2-Chloronaphthalene	1.295	1.264	2.4	111	0.00
48		2-Nitroaniline	0.380	0.385	-1.3	111	0.00
49		Acenaphthylene	1.982	1.943	2.0	112	0.00
50		Dimethylphthalate	1.617	1.566	3.2	112	0.00
51		2,6-Dinitrotoluene	0.367	0.362	1.4	112	0.00
52	C	Acenaphthene	1.201	1.170	2.6	112	0.00
53		3-Nitroaniline	0.392	0.393	-0.3	111	0.00
54	P	2,4-Dinitrophenol	0.180	0.170	5.6	105	0.00
55		Dibenzofuran	1.954	1.873	4.1	111	0.00
56	P	4-Nitrophenol	0.247	0.245	0.8	109	0.00
57		2,4-Dinitrotoluene	0.498	0.500	-0.4	112	0.00
58		Fluorene	1.508	1.455	3.5	111	0.00
59		2,3,4,6-Tetrachlorophenol	0.383	0.385	-0.5	114	0.00
60		Diethylphthalate	1.641	1.612	1.8	113	0.00
61		4-Chlorophenyl-phenylether	0.838	0.817	2.5	113	0.00
62		4-Nitroaniline	0.383	0.393	-2.6	110	0.00
63		Azobenzene	1.517	1.489	1.8	113	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	117	0.00
65		4,6-Dinitro-2-methylphenol	0.140	0.135	3.6	111	0.00
66	c	n-Nitrosodiphenylamine	0.626	0.603	3.7	112	0.00
67		4-Bromophenyl-phenylether	0.225	0.220	2.2	112	0.00
68		Hexachlorobenzene	0.257	0.248	3.5	111	0.00
69		Atrazine	0.222	0.218	1.8	111	0.00
70	C	Pentachlorophenol	0.107	0.123	-15.0	125	0.00
71		Phenanthrene	1.051	1.009	4.0	111	0.00
72		Anthracene	1.049	1.016	3.1	111	0.00
73		Carbazole	1.087	1.047	3.7	111	0.00
74		Di-n-butylphthalate	1.267	1.225	3.3	111	0.00
75	C	Fluoranthene	1.239	1.193	3.7	112	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	117	0.00
77		Benzidine	0.586	0.623	-6.3	123	0.00
78		Pyrene	1.175	1.130	3.8	112	0.00
79	S	Terphenyl-d14	0.935	0.892	4.6	110	0.00
80		Butylbenzylphthalate	0.546	0.524	4.0	110	0.00
81		Benzo(a)anthracene	1.165	1.099	5.7	109	0.00
82		3,3'-Dichlorobenzidine	0.482	0.455	5.6	108	0.00
83		Chrysene	1.120	1.070	4.5	111	0.00
84		Bis(2-ethylhexyl)phthalate	0.809	0.773	4.4	111	0.00
85	c	Di-n-octyl phthalate	1.389	1.300	6.4	108	0.00
86		Indeno(1,2,3-cd)pyrene	1.283	1.132	11.8	99	0.00
87	I	Perylene-d12	1.000	1.000	0.0	111	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN102418\
Data File : BN003288.D
Acq On : 25 Oct 2018 00:59
Operator : MJ/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_N
LabSampleId : SSTDCCC040

Quant Time: Oct 25 05:26:26 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\8270-BN102418.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 24 14:21:25 2018
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	1.101	1.043	5.3	105	0.00
89	Benzo(k)fluoranthene	1.080	1.044	3.3	110	0.00
90 C	Benzo(a)pyrene	1.049	1.000	4.7	106	0.00
91	Dibenzo(a,h)anthracene	1.039	0.939	9.6	99	0.00
92	Benzo(g,h,i)perylene	0.997	0.897	10.0	98	0.00

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