Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG121418\
Data File : BG038571.D
Acq On : 14 Dec 2018 17:09
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 14 23:03:02 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 12 15:26:27 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	115	0.00
2		1,4-Dioxane	0.525	0.478	9.0	103	0.00
3		Pyridine	1.445	1.351	6.5	90	0.00
4		n-Nitrosodimethylamine	0.708	0.724	-2.3	108	0.00
5	S	2-Fluorophenol	1.128	1.151	-2.0	118	0.00
6		Aniline	1.976	1.992	-0.8	115	0.00
7	S	Phenol-d6	1.548	1.559	-0.7	117	0.00
8		2-Chlorophenol	1.305	1.323	-1.4	116	0.00
9		Benzaldehyde	1.004	0.951	5.3	118	0.00
10	C	Phenol	1.645	1.645	0.0	115	0.00
11		bis(2-Chloroethyl)ether	1.309	1.262	3.6	113	0.00
12		1,3-Dichlorobenzene	1.543	1.530	0.8	116	0.00
13	C	1,4-Dichlorobenzene	1.580	1.550	1.9	115	0.00
14		1,2-Dichlorobenzene	1.490	1.478	0.8	118	0.00
15		Benzyl Alcohol	1.208	1.269	-5.0	116	0.00
16		2,2'-oxybis(1-Chloropropane	2.999	2.910	3.0	113	0.00
17		2-Methylphenol	1.102	1.126	-2.2	115	0.00
18		Hexachloroethane	0.577	0.575	0.3	117	0.00
19	P	n-Nitroso-di-n-propylamine	1.087	1.106	-1.7	117	0.00
20		3+4-Methylphenols	1.522	1.580	-3.8	117	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	115	0.00
22		Acetophenone	0.500	0.509	-1.8	116	0.00
23	S	Nitrobenzene-d5	0.391	0.406	-3.8	118	0.00
24		Nitrobenzene	0.396	0.412	-4.0	119	0.00
25		Isophorone	0.703	0.674	4.1	93	0.00
26	C	2-Nitrophenol	0.203	0.210	-3.4	118	0.00
27		2,4-Dimethylphenol	0.291	0.299	-2.7	117	0.00
28		bis(2-Chloroethoxy)methane	0.397	0.410	-3.3	116	0.00
29	C	2,4-Dichlorophenol	0.323	0.351	-8.7	119	0.00
30		1,2,4-Trichlorobenzene	0.390	0.400	-2.6	118	0.00
31		Naphthalene	0.985	0.993	-0.8	116	0.00
32		Benzoic acid	0.154	0.190	-23.4	128	0.00
33		4-Chloroaniline	0.428	0.456	-6.5	118	0.00
34	C	Hexachlorobutadiene	0.264	0.271	-2.7	115	0.00
35		Caprolactam	0.134	0.135	-0.7	121	0.00
36	C	4-Chloro-3-methylphenol	0.369	0.375	-1.6	117	0.00
37		2-Methylnaphthalene	0.703	0.717	-2.0	117	0.00
38		1-Methylnaphthalene	0.682	0.693	-1.6	117	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
40		1,2,4,5-Tetrachlorobenzene	0.748	0.751	-0.4	115	0.00
41	P	Hexachlorocyclopentadiene	0.411	0.389	5.4	107	0.00
42	S	2,4,6-Tribromophenol	0.276	0.292	-5.8	119	0.00
43	C	2,4,6-Trichlorophenol	0.460	0.490	-6.5	116	0.00
44		2,4,5-Trichlorophenol	0.487	0.524	-7.6	119	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG121418\
Data File : BG038571.D
Acq On : 14 Dec 2018 17:09
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 14 23:03:02 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 12 15:26:27 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	1.458	1.467	-0.6	116	0.00
46		1,1'-Biphenyl	1.677	1.696	-1.1	115	0.00
47		2-Chloronaphthalene	1.295	1.289	0.5	115	0.00
48		2-Nitroaniline	0.413	0.444	-7.5	121	0.00
49		Acenaphthylene	1.936	1.964	-1.4	116	0.00
50		Dimethylphthalate	1.633	1.685	-3.2	118	0.00
51		2,6-Dinitrotoluene	0.370	0.384	-3.8	119	0.00
52	C	Acenaphthene	1.158	1.167	-0.8	116	0.00
53		3-Nitroaniline	0.377	0.403	-6.9	120	0.00
54	P	2,4-Dinitrophenol	0.193	0.173	10.4	100	0.00
55		Dibenzofuran	1.985	1.986	-0.1	117	0.00
56	P	4-Nitrophenol	0.286	0.298	-4.2	114	0.00
57		2,4-Dinitrotoluene	0.521	0.540	-3.6	116	0.00
58		Fluorene	1.470	1.496	-1.8	116	0.00
59		2,3,4,6-Tetrachlorophenol	0.438	0.453	-3.4	114	0.00
60		Diethylphthalate	1.793	1.811	-1.0	118	0.00
61		4-Chlorophenyl-phenylether	0.917	0.938	-2.3	117	0.00
62		4-Nitroaniline	0.388	0.424	-9.3	120	0.00
63		Azobenzene	1.498	1.527	-1.9	116	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	117	0.00
65		4,6-Dinitro-2-methylphenol	0.143	0.136	4.9	109	0.00
66	c	n-Nitrosodiphenylamine	0.588	0.592	-0.7	118	0.00
67		4-Bromophenyl-phenylether	0.244	0.246	-0.8	117	0.00
68		Hexachlorobenzene	0.253	0.255	-0.8	115	0.00
69		Atrazine	0.218	0.227	-4.1	118	0.00
70	C	Pentachlorophenol	0.150	0.154	-2.7	117	0.00
71		Phenanthrene	1.037	1.037	0.0	118	0.00
72		Anthracene	1.024	1.030	-0.6	116	0.00
73		Carbazole	1.013	1.030	-1.7	118	0.00
74		Di-n-butylphthalate	1.162	1.184	-1.9	117	0.00
75	C	Fluoranthene	1.283	1.257	2.0	117	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	112	0.00
77		Benzidine	0.591	0.639	-8.1	101	0.00
78		Pyrene	1.321	1.336	-1.1	116	0.00
79	S	Terphenyl-d14	0.919	0.949	-3.3	116	0.00
80		Butylbenzylphthalate	0.516	0.547	-6.0	116	0.00
81		Benzo(a)anthracene	1.219	1.243	-2.0	112	0.00
82		3,3'-Dichlorobenzidine	0.483	0.514	-6.4	113	0.00
83		Chrysene	1.174	1.182	-0.7	112	0.00
84		Bis(2-ethylhexyl)phthalate	0.732	0.777	-6.1	114	0.00
85	c	Di-n-octyl phthalate	1.226	1.303	-6.3	113	0.00
86		Indeno(1,2,3-cd)pyrene	1.380	1.364	1.2	106	0.00
87	I	Perylene-d12	1.000	1.000	0.0	110	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA G\Data\BG121418\
Data File : BG038571.D
Acq On : 14 Dec 2018 17:09
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 14 23:03:02 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA G\METHODS\8270-BG121218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 12 15:26:27 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	1.098	1.126	-2.6	110	0.00
89	Benzo(k)fluoranthene	1.093	1.095	-0.2	108	0.00
90 C	Benzo(a)pyrene	1.059	1.086	-2.5	110	0.00
91	Dibenzo(a,h)anthracene	1.055	1.042	1.2	106	0.00
92	Benzo(g,h,i)perylene	1.039	1.013	2.5	105	0.00

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