Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081018\
Data File : BG036287.D
Acq On : 11 Aug 2018 4:35
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 11 05:09:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 31 12:30:06 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	67	-0.02
2		1,4-Dioxane	0.613	0.523	14.7	62	-0.01
3		Pyridine	1.601	1.334	16.7	54	0.00
4		n-Nitrosodimethylamine	0.687	0.562	18.2	55	-0.01
5	S	2-Fluorophenol	1.205	1.087	9.8	60	0.00
6		Aniline	2.330	1.934	17.0	56	-0.02
7	S	Phenol-d6	1.797	1.580	12.1	58	-0.01
8		2-Chlorophenol	1.225	1.143	6.7	62	-0.02
9		Benzaldehyde	1.103	0.895	18.9	53	-0.01
10	C	Phenol	1.816	1.619	10.8	59	-0.01
11		bis(2-Chloroethyl)ether	1.422	1.194	16.0	56	-0.02
12		1,3-Dichlorobenzene	1.480	1.405	5.1	65	-0.03
13	C	1,4-Dichlorobenzene	1.503	1.440	4.2	65	-0.03
14		1,2-Dichlorobenzene	1.444	1.376	4.7	65	-0.03
15		Benzyl Alcohol	1.632	1.321	19.1	54	-0.01
16		2,2'-oxybis(1-Chloropropane	1.192	0.867	27.3#	49#	-0.03
17		2-Methylphenol	1.235	1.080	12.6	57	-0.01
18		Hexachloroethane	0.575	0.540	6.1	65	-0.02
19	P	n-Nitroso-di-n-propylamine	1.513	1.179	22.1	53	-0.03
20		3+4-Methylphenols	1.713	1.460	14.8	55	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	58	-0.03
22		Acetophenone	0.622	0.574	7.7	56	-0.02
23	S	Nitrobenzene-d5	0.479	0.458	4.4	57	-0.02
24		Nitrobenzene	0.481	0.461	4.2	58	-0.02
25		Isophorone	0.889	0.760	14.5	51	-0.02
26	C	2-Nitrophenol	0.171	0.184	-7.6	62	-0.02
27		2,4-Dimethylphenol	0.289	0.267	7.6	54	-0.02
28		bis(2-Chloroethoxy)methane	0.490	0.433	11.6	53	-0.03
29	C	2,4-Dichlorophenol	0.330	0.343	-3.9	60	-0.01
30		1,2,4-Trichlorobenzene	0.405	0.418	-3.2	62	-0.02
31		Naphthalene	1.042	0.987	5.3	58	-0.02
32		Benzoic acid	0.195	0.089	54.4#	27#	0.00
33		4-Chloroaniline	0.454	0.425	6.4	57	-0.01
34	C	Hexachlorobutadiene	0.316	0.347	-9.8	66	-0.03
35		Caprolactam	0.142	0.121	14.8	53	-0.01
36	C	4-Chloro-3-methylphenol	0.443	0.438	1.1	58	-0.01
37		2-Methylnaphthalene	0.776	0.758	2.3	58	-0.03
38	I	Acenaphthene-d10	1.000	1.000	0.0	58	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.755	0.765	-1.3	63	-0.02
40	P	Hexachlorocyclopentadiene	0.396	0.470	-18.7	71	-0.02
41	S	2,4,6-Tribromophenol	0.264	0.298	-12.9	69	-0.02
42	C	2,4,6-Trichlorophenol	0.441	0.465	-5.4	61	-0.02
43		2,4,5-Trichlorophenol	0.494	0.514	-4.0	65	0.00
44	S	2-Fluorobiphenyl	1.493	1.467	1.7	61	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081018\
Data File : BG036287.D
Acq On : 11 Aug 2018 4:35
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 11 05:09:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 31 12:30:06 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		1,1'-Biphenyl	1.551	1.498	3.4	59	-0.03
46		2-Chloronaphthalene	1.206	1.140	5.5	59	-0.02
47		2-Nitroaniline	0.426	0.412	3.3	58	-0.01
48		Acenaphthylene	2.020	1.930	4.5	59	-0.02
49		Dimethylphthalate	1.752	1.634	6.7	59	-0.02
50		2,6-Dinitrotoluene	0.357	0.385	-7.8	65	-0.02
51	C	Acenaphthene	1.259	1.206	4.2	60	-0.02
52		3-Nitroaniline	0.365	0.355	2.7	58	-0.02
53	P	2,4-Dinitrophenol	0.158	0.135	14.6	52	0.00
54		Dibenzofuran	2.002	1.961	2.0	60	-0.02
55	P	4-Nitrophenol	0.260	0.239	8.1	55	0.00
56		2,4-Dinitrotoluene	0.512	0.557	-8.8	63	-0.01
57		Fluorene	1.678	1.652	1.5	62	-0.02
58		2,3,4,6-Tetrachlorophenol	0.473	0.496	-4.9	63	-0.02
59		Diethylphthalate	1.802	1.715	4.8	59	-0.02
60		4-Chlorophenyl-phenylether	0.986	1.005	-1.9	64	-0.03
61		4-Nitroaniline	0.382	0.360	5.8	56	-0.01
62		Azobenzene	1.777	1.620	8.8	56	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	62	-0.02
64		4,6-Dinitro-2-methylphenol	0.111	0.104	6.3	60	-0.01
65	c	n-Nitrosodiphenylamine	0.569	0.554	2.6	63	-0.02
66		4-Bromophenyl-phenylether	0.232	0.233	-0.4	64	-0.02
67		Hexachlorobenzene	0.239	0.242	-1.3	66	-0.02
68		Atrazine	0.234	0.217	7.3	58	-0.01
69	C	Pentachlorophenol	0.146	0.143	2.1	61	-0.01
70		Phenanthrene	0.998	0.980	1.8	64	-0.03
71		Anthracene	0.994	0.972	2.2	64	-0.02
72		Carbazole	0.944	0.905	4.1	62	-0.01
73		Di-n-butylphthalate	1.115	1.086	2.6	63	-0.02
74	C	Fluoranthene	1.344	1.333	0.8	64	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	64	-0.02
76		Benzidine	0.526	0.275	47.7#	36#	-0.01
77		Pyrene	1.265	1.207	4.6	65	-0.02
78	S	Terphenyl-d14	0.907	0.918	-1.2	68	-0.02
79		Butylbenzylphthalate	0.452	0.440	2.7	63	-0.01
80		Benzo(a)anthracene	1.246	1.196	4.0	65	-0.02
81		3,3'-Dichlorobenzidine	0.435	0.410	5.7	62	-0.03
82		Chrysene	1.155	1.120	3.0	66	-0.02
83		Bis(2-ethylhexyl)phthalate	0.615	0.616	-0.2	65	-0.03
84	c	Di-n-octyl phthalate	1.029	1.035	-0.6	65	-0.03
85		Indeno(1,2,3-cd)pyrene	1.279	1.201	6.1	62	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	65	-0.04
87		Benzo(b)fluoranthene	1.216	1.196	1.6	67	-0.03

Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081018\
Data File : BG036287.D
Acq On : 11 Aug 2018 4:35
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 11 05:09:53 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_G\METHODS\8270-BG071218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 31 12:30:06 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(k)fluoranthene	1.188	1.151	3.1	64	-0.03
89 C	Benzo(a)pyrene	1.131	1.113	1.6	65	-0.03
90	Dibenzo(a,h)anthracene	1.110	0.999	10.0	60	-0.06
91	Benzo(g,h,i)perylene	1.086	0.995	8.4	60	-0.05

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