Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071816\
Data File : BG023157.D
Acq On : 18 Jul 2016 11:36
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 19 01:01:02 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 18:30:58 2016
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	109	0.00
2		1,4-Dioxane	0.475	0.447	5.9	107	0.00
3		Pyridine	1.627	1.651	-1.5	111	0.00
4		n-Nitrosodimethylamine	0.698	0.780	-11.7	121	0.00
5	S	2-Fluorophenol	1.223	1.202	1.7	110	0.00
6		Aniline	2.654	2.745	-3.4	112	0.00
7	S	Phenol-d6	1.915	1.945	-1.6	109	0.00
8		2-Chlorophenol	1.301	1.401	-7.7	118	0.00
9		Benzaldehyde	1.280	1.083	15.4	94	0.00
10	C	Phenol	1.971	2.124	-7.8	117	0.00
11		bis(2-Chloroethyl)ether	1.562	1.596	-2.2	112	0.00
12		1,3-Dichlorobenzene	1.593	1.536	3.6	109	0.00
13	C	1,4-Dichlorobenzene	1.595	1.605	-0.6	112	0.00
14		1,2-Dichlorobenzene	1.585	1.547	2.4	111	0.00
15		Benzyl Alcohol	1.754	2.007	-14.4	122	0.00
16		2,2'-oxybis(1-Chloropropane	1.908	1.965	-3.0	113	0.00
17		2-Methylphenol	1.283	1.416	-10.4	120	0.00
18		Hexachloroethane	0.673	0.649	3.6	114	0.00
19	P	n-Nitroso-di-n-propylamine	1.727	1.822	-5.5	112	0.00
20		3+4-Methylphenols	1.789	2.031	-13.5	117	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	113	0.00
22		Acetophenone	0.600	0.577	3.8	109	0.00
23	S	Nitrobenzene-d5	0.467	0.465	0.4	113	0.00
24		Nitrobenzene	0.496	0.509	-2.6	119	0.00
25		Isophorone	0.939	0.953	-1.5	111	0.00
26	C	2-Nitrophenol	0.195	0.194	0.5	106	0.00
27		2,4-Dimethylphenol	0.337	0.328	2.7	111	0.00
28		bis(2-Chloroethoxy)methane	0.524	0.500	4.6	106	0.00
29	C	2,4-Dichlorophenol	0.359	0.361	-0.6	111	0.00
30		1,2,4-Trichlorobenzene	0.411	0.377	8.3	104	0.00
31		Naphthalene	1.018	0.971	4.6	109	0.00
32		Benzoic acid	0.306	0.225	26.5#	78	0.00
33		4-Chloroaniline	0.498	0.498	0.0	113	0.00
34	C	Hexachlorobutadiene	0.334	0.291	12.9	101	0.00
35		Caprolactam	0.148	0.139	6.1	106	0.00
36	C	4-Chloro-3-methylphenol	0.491	0.470	4.3	106	0.00
37		2-Methylnaphthalene	0.805	0.800	0.6	110	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
39		1,2,4,5-Tetrachlorobenzene	0.862	0.652	24.4	83	0.00
40	P	Hexachlorocyclopentadiene	0.496	0.420	15.3	92	0.00
41	S	2,4,6-Tribromophenol	0.359	0.288	19.8	89	0.00
42	C	2,4,6-Trichlorophenol	0.495	0.472	4.6	102	0.00
43		2,4,5-Trichlorophenol	0.509	0.489	3.9	106	0.00
44	S	2-Fluorobiphenyl	1.270	1.177	7.3	103	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071816\
Data File : BG023157.D
Acq On : 18 Jul 2016 11:36
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 19 01:01:02 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 18:30:58 2016
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	1.501	1.415	5.7	102	0.00
46		2-Chloronaphthalene	1.217	1.184	2.7	105	0.00
47		2-Nitroaniline	0.519	0.515	0.8	108	0.00
48		Acenaphthylene	1.826	1.774	2.8	106	0.00
49		Dimethylphthalate	1.634	1.546	5.4	105	0.00
50		2,6-Dinitrotoluene	0.367	0.355	3.3	106	0.00
51	C	Acenaphthene	1.193	1.174	1.6	108	0.00
52		3-Nitroaniline	0.366	0.372	-1.6	110	0.00
53	P	2,4-Dinitrophenol	0.252	0.207	17.9	86	0.00
54		Dibenzofuran	1.786	1.828	-2.4	114	0.00
55	P	4-Nitrophenol	0.307	0.286	6.8	101	0.00
56		2,4-Dinitrotoluene	0.535	0.526	1.7	108	0.00
57		Fluorene	1.537	1.494	2.8	108	0.00
58		2,3,4,6-Tetrachlorophenol	0.509	0.488	4.1	104	0.00
59		Diethylphthalate	1.662	1.598	3.9	106	0.00
60		4-Chlorophenyl-phenylether	0.936	0.857	8.4	99	0.00
61		4-Nitroaniline	0.396	0.378	4.5	105	0.00
62		Azobenzene	1.590	1.784	-12.2	124	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	102	0.00
64		4,6-Dinitro-2-methylphenol	0.149	0.137	8.1	90	0.00
65	c	n-Nitrosodiphenylamine	0.576	0.588	-2.1	103	0.00
66		4-Bromophenyl-phenylether	0.260	0.245	5.8	95	0.00
67		Hexachlorobenzene	0.283	0.263	7.1	94	0.00
68		Atrazine	0.256	0.246	3.9	99	0.00
69	C	Pentachlorophenol	0.183	0.160	12.6	85	0.00
70		Phenanthrene	0.980	0.996	-1.6	105	0.00
71		Anthracene	0.992	1.007	-1.5	105	0.00
72		Carbazole	0.858	0.879	-2.4	106	0.00
73		Di-n-butylphthalate	0.954	1.026	-7.5	107	0.00
74	C	Fluoranthene	1.203	1.103	8.3	99	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	96	0.00
76		Benzidine	0.573	0.447	22.0	73	0.00
77		Pyrene	1.124	1.075	4.4	98	0.00
78	S	Terphenyl-d14	0.646	0.722	-11.8	108	0.00
79		Butylbenzylphthalate	0.445	0.487	-9.4	106	0.00
80		Benzo(a)anthracene	1.119	1.108	1.0	98	0.00
81		3,3'-Dichlorobenzidine	0.491	0.446	9.2	86	0.00
82		Chrysene	1.050	1.025	2.4	97	0.00
83		Bis(2-ethylhexyl)phthalate	0.618	0.664	-7.4	106	0.00
84	c	Di-n-octyl phthalate	1.031	1.084	-5.1	102	0.00
85		Indeno(1,2,3-cd)pyrene	1.338	1.208	9.7	90	0.00
86	I	Perylene-d12	1.000	1.000	0.0	90	0.00
87		Benzo(b)fluoranthene	1.154	1.201	-4.1	96	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG071816\
Data File : BG023157.D
Acq On : 18 Jul 2016 11:36
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 19 01:01:02 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 18:30:58 2016
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	1.095	1.128	-3.0	92	0.00
89 C	Benzo(a)pyrene	1.106	1.152	-4.2	95	0.00
90	Dibenzo(a,h)anthracene	1.098	1.030	6.2	86	0.00
91	Benzo(g,h,i)perylene	1.099	1.006	8.5	83	0.00

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