Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062716\
Data File : BG022634.D
Acq On : 27 Jun 2016 9:02
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 27 14:49:58 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 24 20:14:59 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	129	0.00
2		1,4-Dioxane	0.505	0.492	2.6	123	0.00
3		Pyridine	1.414	1.629	-15.2	145	-0.02
4		n-Nitrosodimethylamine	0.809	0.722	10.8	112	-0.01
5	S	2-Fluorophenol	1.247	1.217	2.4	126	0.00
6		Aniline	2.330	2.510	-7.7	136	0.00
7	S	Phenol-d6	1.758	1.793	-2.0	131	0.00
8		2-Chlorophenol	1.292	1.280	0.9	131	0.00
9		Benzaldehyde	0.619	0.762	-23.1	156#	0.00
10	C	Phenol	1.858	1.942	-4.5	133	0.00
11		bis(2-Chloroethyl)ether	1.529	1.444	5.6	115	0.00
12		1,3-Dichlorobenzene	1.572	1.530	2.7	125	0.00
13	C	1,4-Dichlorobenzene	1.576	1.538	2.4	128	0.00
14		1,2-Dichlorobenzene	1.498	1.463	2.3	129	0.00
15		Benzyl Alcohol	1.626	1.747	-7.4	129	0.00
16		2,2'-oxybis(1-Chloropropane	1.691	1.609	4.8	122	0.00
17		2-Methylphenol	1.224	1.261	-3.0	136	0.00
18		Hexachloroethane	0.652	0.642	1.5	124	-0.01
19	P	n-Nitroso-di-n-propylamine	1.464	1.447	1.2	127	0.00
20		3+4-Methylphenols	1.687	1.739	-3.1	135	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	133	0.00
22		Acetophenone	0.578	0.559	3.3	126	0.00
23	S	Nitrobenzene-d5	0.473	0.461	2.5	128	0.00
24		Nitrobenzene	0.522	0.507	2.9	131	0.00
25		Isophorone	0.917	0.882	3.8	129	0.00
26	C	2-Nitrophenol	0.188	0.196	-4.3	140	0.00
27		2,4-Dimethylphenol	0.359	0.320	10.9	124	0.00
28		bis(2-Chloroethoxy)methane	0.501	0.488	2.6	133	0.00
29	C	2,4-Dichlorophenol	0.350	0.361	-3.1	135	0.00
30		1,2,4-Trichlorobenzene	0.416	0.403	3.1	130	0.00
31		Naphthalene	1.005	0.953	5.2	127	0.00
32		Benzoic acid	0.216	0.260	-20.4	154#	0.08
33		4-Chloroaniline	0.433	0.466	-7.6	135	0.00
34	C	Hexachlorobutadiene	0.323	0.315	2.5	127	0.00
35		Caprolactam	0.110	0.125	-13.6	150	0.03
36	C	4-Chloro-3-methylphenol	0.430	0.446	-3.7	138	0.00
37		2-Methylnaphthalene	0.780	0.752	3.6	129	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	141	0.00
39		1,2,4,5-Tetrachlorobenzene	0.755	0.724	4.1	139	0.00
40	P	Hexachlorocyclopentadiene	0.603	0.523	13.3	119	0.00
41	S	2,4,6-Tribromophenol	0.315	0.328	-4.1	146	0.00
42	C	2,4,6-Trichlorophenol	0.484	0.490	-1.2	139	0.00
43		2,4,5-Trichlorophenol	0.507	0.528	-4.1	149	0.00
44	S	2-Fluorobiphenyl	1.261	1.202	4.7	128	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062716\
Data File : BG022634.D
Acq On : 27 Jun 2016 9:02
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 27 14:49:58 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 24 20:14:59 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	1.524	1.410	7.5	131	0.00
46		2-Chloronaphthalene	1.254	1.177	6.1	133	0.00
47		2-Nitroaniline	0.484	0.488	-0.8	140	0.00
48		Acenaphthylene	1.818	1.722	5.3	133	0.00
49		Dimethylphthalate	1.659	1.561	5.9	134	0.00
50		2,6-Dinitrotoluene	0.361	0.365	-1.1	141	0.00
51	C	Acenaphthene	1.339	1.288	3.8	133	0.00
52		3-Nitroaniline	0.337	0.355	-5.3	146	0.00
53	P	2,4-Dinitrophenol	0.222	0.216	2.7	133	0.00
54		Dibenzofuran	1.940	1.807	6.9	133	0.00
55	P	4-Nitrophenol	0.285	0.309	-8.4	155#	0.02
56		2,4-Dinitrotoluene	0.499	0.529	-6.0	148	0.00
57		Fluorene	1.548	1.489	3.8	136	0.00
58		2,3,4,6-Tetrachlorophenol	0.525	0.557	-6.1	150#	0.00
59		Diethylphthalate	1.706	1.592	6.7	135	0.00
60		4-Chlorophenyl-phenylether	0.922	0.891	3.4	138	0.00
61		4-Nitroaniline	0.348	0.366	-5.2	147	0.02
62		Azobenzene	1.847	1.692	8.4	129	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	149	0.00
64		4,6-Dinitro-2-methylphenol	0.134	0.138	-3.0	143	0.00
65	c	n-Nitrosodiphenylamine	0.582	0.552	5.2	138	0.00
66		4-Bromophenyl-phenylether	0.259	0.246	5.0	141	0.00
67		Hexachlorobenzene	0.274	0.262	4.4	143	0.00
68		Atrazine	0.246	0.248	-0.8	148	0.00
69	C	Pentachlorophenol	0.189	0.184	2.6	144	0.00
70		Phenanthrene	1.020	0.937	8.1	136	0.00
71		Anthracene	1.050	0.959	8.7	135	0.00
72		Carbazole	0.890	0.846	4.9	139	0.00
73		Di-n-butylphthalate	1.036	0.968	6.6	131	0.00
74	C	Fluoranthene	1.175	1.091	7.1	133	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	147	0.00
76		Benzidine	0.213	0.390	-83.1#	275#	-0.02
77		Pyrene	1.065	1.026	3.7	133	0.00
78	S	Terphenyl-d14	0.755	0.609	19.3	125	0.00
79		Butylbenzylphthalate	0.461	0.455	1.3	138	0.00
80		Benzo(a)anthracene	1.107	1.092	1.4	137	0.00
81		3,3'-Dichlorobenzidine	0.457	0.451	1.3	139	0.00
82		Chrysene	1.040	1.016	2.3	136	0.00
83		Bis(2-ethylhexyl)phthalate	0.649	0.605	6.8	134	0.00
84	c	Di-n-octyl phthalate	1.070	0.985	7.9	132	0.00
85		Indeno(1,2,3-cd)pyrene	1.367	1.363	0.3	145	0.03
86	I	Perylene-d12	1.000	1.000	0.0	146	0.02
87		Benzo(b)fluoranthene	1.203	1.153	4.2	139	0.02

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG062716\
Data File : BG022634.D
Acq On : 27 Jun 2016 9:02
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jun 27 14:49:58 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG062516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jun 24 20:14:59 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	1.137	1.077	5.3	139	0.02
89 C	Benzo(a)pyrene	1.172	1.116	4.8	139	0.02
90	Dibenzo(a,h)anthracene	1.104	1.079	2.3	146	0.04
91	Benzo(g,h,i)perylene	1.123	1.086	3.3	144	0.04

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