Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061417\
Data File : BG027423.D
Acq On : 14 Jun 2017 18:48
Operator : SJ/MA
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG061417

Quant Time: Jun 14 19:22:50 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG061417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 14 19:09:08 2017
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	102	0.00
2		1,4-Dioxane	0.586	0.604	-3.1	106	0.00
3		Pyridine	1.792	1.894	-5.7	104	0.00
4		n-Nitrosodimethylamine	0.773	0.763	1.3	101	0.00
5	S	2-Fluorophenol	1.415	1.447	-2.3	100	0.00
6		Aniline	2.589	2.729	-5.4	102	0.00
7	S	Phenol-d6	2.175	2.230	-2.5	100	0.00
8		2-Chlorophenol	1.477	1.526	-3.3	102	0.00
9		Benzaldehyde	1.465	1.514	-3.3	101	0.00
10	C	Phenol	2.242	2.305	-2.8	101	0.00
11		bis(2-Chloroethyl)ether	1.778	1.803	-1.4	102	0.00
12		1,3-Dichlorobenzene	1.554	1.597	-2.8	102	0.00
13	C	1,4-Dichlorobenzene	1.615	1.627	-0.7	101	0.00
14		1,2-Dichlorobenzene	1.586	1.580	0.4	100	0.00
15		Benzyl Alcohol	1.575	1.627	-3.3	101	0.00
16		2,2'-oxybis(1-Chloropropane	3.068	3.049	0.6	99	0.00
17		2-Methylphenol	1.515	1.561	-3.0	101	0.00
18		Hexachloroethane	0.594	0.588	1.0	102	0.00
19	P	n-Nitroso-di-n-propylamine	1.645	1.726	-4.9	101	0.00
20		3+4-Methylphenols	2.180	2.227	-2.2	99	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.537	0.547	-1.9	99	0.00
23	S	Nitrobenzene-d5	0.392	0.413	-5.4	101	0.00
24		Nitrobenzene	0.434	0.453	-4.4	100	0.00
25		Isophorone	0.858	0.902	-5.1	99	0.00
26	C	2-Nitrophenol	0.169	0.175	-3.6	101	0.00
27		2,4-Dimethylphenol	0.347	0.345	0.6	98	0.00
28		bis(2-Chloroethoxy)methane	0.523	0.530	-1.3	97	0.00
29	C	2,4-Dichlorophenol	0.300	0.311	-3.7	98	0.00
30		1,2,4-Trichlorobenzene	0.311	0.324	-4.2	103	0.00
31		Naphthalene	1.092	1.101	-0.8	99	0.00
32		Benzoic acid	0.199	0.220	-10.6	104	0.00
33		4-Chloroaniline	0.460	0.496	-7.8	101	0.00
34	C	Hexachlorobutadiene	0.191	0.199	-4.2	102	0.00
35		Caprolactam	0.133	0.144	-8.3	101	0.00
36	C	4-Chloro-3-methylphenol	0.429	0.452	-5.4	100	0.00
37		2-Methylnaphthalene	0.795	0.820	-3.1	99	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
39		1,2,4,5-Tetrachlorobenzene	0.589	0.600	-1.9	101	0.00
40	P	Hexachlorocyclopentadiene	0.328	0.326	0.6	100	0.00
41	S	2,4,6-Tribromophenol	0.300	0.314	-4.7	98	0.00
42	C	2,4,6-Trichlorophenol	0.401	0.412	-2.7	97	0.00
43		2,4,5-Trichlorophenol	0.456	0.473	-3.7	99	0.00
44	S	2-Fluorobiphenyl	1.384	1.402	-1.3	98	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061417\
Data File : BG027423.D
Acq On : 14 Jun 2017 18:48
Operator : SJ/MA
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG061417

Quant Time: Jun 14 19:22:50 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG061417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 14 19:09:08 2017
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	1.591	1.594	-0.2	97	0.00
46		2-Chloronaphthalene	1.204	1.213	-0.7	98	0.00
47		2-Nitroaniline	0.476	0.506	-6.3	99	0.00
48		Acenaphthylene	2.103	2.155	-2.5	99	0.00
49		Dimethylphthalate	1.661	1.665	-0.2	97	0.00
50		2,6-Dinitrotoluene	0.342	0.371	-8.5	101	0.00
51	C	Acenaphthene	1.402	1.435	-2.4	100	0.00
52		3-Nitroaniline	0.378	0.390	-3.2	100	0.00
53	P	2,4-Dinitrophenol	0.186	0.198	-6.5	102	0.00
54		Dibenzofuran	1.911	1.927	-0.8	99	0.00
55	P	4-Nitrophenol	0.299	0.320	-7.0	100	0.00
56		2,4-Dinitrotoluene	0.488	0.513	-5.1	102	0.00
57		Fluorene	1.718	1.752	-2.0	99	0.00
58		2,3,4,6-Tetrachlorophenol	0.398	0.415	-4.3	100	0.00
59		Diethylphthalate	1.707	1.731	-1.4	99	0.00
60		4-Chlorophenyl-phenylether	0.866	0.870	-0.5	97	0.00
61		4-Nitroaniline	0.387	0.414	-7.0	100	0.00
62		Azobenzene	1.757	1.795	-2.2	98	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
64		4,6-Dinitro-2-methylphenol	0.116	0.125	-7.8	105	0.00
65	c	n-Nitrosodiphenylamine	0.624	0.645	-3.4	99	0.00
66		4-Bromophenyl-phenylether	0.238	0.242	-1.7	97	0.00
67		Hexachlorobenzene	0.263	0.269	-2.3	99	0.00
68		Atrazine	0.217	0.227	-4.6	100	0.00
69	C	Pentachlorophenol	0.151	0.160	-6.0	101	0.00
70		Phenanthrene	1.116	1.142	-2.3	99	0.00
71		Anthracene	1.129	1.162	-2.9	99	0.00
72		Carbazole	1.081	1.116	-3.2	99	0.00
73		Di-n-butylphthalate	1.179	1.231	-4.4	100	0.00
74	C	Fluoranthene	1.285	1.324	-3.0	100	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	99	0.00
76		Benzidine	0.439	0.466	-6.2	102	0.00
77		Pyrene	1.103	1.139	-3.3	100	0.00
78	S	Terphenyl-d14	0.780	0.807	-3.5	99	0.00
79		Butylbenzylphthalate	0.459	0.479	-4.4	100	0.00
80		Benzo(a)anthracene	1.150	1.185	-3.0	101	0.00
81		3,3'-Dichlorobenzidine	0.436	0.463	-6.2	100	0.00
82		Chrysene	1.099	1.118	-1.7	99	0.00
83		Bis(2-ethylhexyl)phthalate	0.665	0.708	-6.5	101	0.00
84	c	Di-n-octyl phthalate	1.114	1.179	-5.8	101	0.00
85		Indeno(1,2,3-cd)pyrene	1.360	1.390	-2.2	99	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	100	0.00
87		Benzo(b)fluoranthene	1.267	1.304	-2.9	99	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG061417\
Data File : BG027423.D
Acq On : 14 Jun 2017 18:48
Operator : SJ/MA
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG061417

Quant Time: Jun 14 19:22:50 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG061417.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 14 19:09:08 2017
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	1.247	1.268	-1.7	99	0.00
89 C	Benzo(a)pyrene	1.188	1.215	-2.3	99	0.00
90	Dibenzo(a,h)anthracene	1.278	1.304	-2.0	98	0.00
91	Benzo(g,h,i)perylene	1.219	1.245	-2.1	99	0.00

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