Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042315\
Data File : BG016659.D
Acq On : 23 Apr 2015 19:09
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG042315

Quant Time: Apr 24 02:19:54 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 23 19:19:11 2015
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	96	0.00
2		1,4-Dioxane	0.534	0.549	-2.8	92	0.00
3		Pyridine	1.516	1.567	-3.4	93	0.00
4		n-Nitrosodimethylamine	0.447	0.451	-0.9	94	0.00
5	S	2-Fluorophenol	1.232	1.249	-1.4	92	0.00
6		Aniline	2.371	2.489	-5.0	93	0.00
7	S	Phenol-d6	1.729	1.803	-4.3	93	0.00
8		2-Chlorophenol	1.523	1.542	-1.2	92	0.00
9		Benzaldehyde	0.706	0.633	10.3	91	0.00
10	C	Phenol	1.823	1.886	-3.5	94	0.00
11		bis(2-Chloroethyl)ether	1.423	1.435	-0.8	93	0.00
12		1,3-Dichlorobenzene	1.575	1.572	0.2	92	0.00
13	C	1,4-Dichlorobenzene	1.610	1.603	0.4	91	0.00
14		1,2-Dichlorobenzene	1.537	1.529	0.5	92	0.00
15		Benzyl Alcohol	1.085	1.154	-6.4	92	0.00
16		2,2'-oxybis(1-Chloropropane	1.332	1.341	-0.7	92	0.00
17		2-Methylphenol	1.219	1.268	-4.0	92	0.00
18		Hexachloroethane	0.568	0.562	1.1	92	0.00
19	P	n-Nitroso-di-n-propylamine	1.078	1.150	-6.7	93	0.00
20		3+4-Methylphenols	1.713	1.809	-5.6	94	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	94	0.00
22		Acetophenone	0.439	0.464	-5.7	95	0.00
23	S	Nitrobenzene-d5	0.317	0.333	-5.0	93	0.00
24		Nitrobenzene	0.332	0.352	-6.0	95	0.00
25		Isophorone	0.656	0.703	-7.2	93	0.00
26	C	2-Nitrophenol	0.193	0.208	-7.8	94	0.00
27		2,4-Dimethylphenol	0.322	0.343	-6.5	95	0.00
28		bis(2-Chloroethoxy)methane	0.396	0.418	-5.6	95	0.00
29	C	2,4-Dichlorophenol	0.274	0.297	-8.4	94	0.00
30		1,2,4-Trichlorobenzene	0.290	0.298	-2.8	93	0.00
31		Naphthalene	1.055	1.100	-4.3	95	0.00
32		Benzoic acid	0.151	0.145	4.0	82	0.00
33		4-Chloroaniline	0.486	0.523	-7.6	95	0.00
34	C	Hexachlorobutadiene	0.130	0.132	-1.5	94	0.00
35		Caprolactam	0.148	0.169	-14.2	94	0.00
36	C	4-Chloro-3-methylphenol	0.318	0.349	-9.7	95	0.00
37		2-Methylnaphthalene	0.762	0.805	-5.6	95	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
39		1,2,4,5-Tetrachlorobenzene	0.476	0.488	-2.5	95	0.00
40	P	Hexachlorocyclopentadiene	0.153	0.148	3.3	88	0.00
41	S	2,4,6-Tribromophenol	0.148	0.158	-6.8	96	0.00
42	C	2,4,6-Trichlorophenol	0.348	0.370	-6.3	94	0.00
43		2,4,5-Trichlorophenol	0.368	0.386	-4.9	94	0.00
44	S	2-Fluorobiphenyl	1.245	1.269	-1.9	95	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042315\
Data File : BG016659.D
Acq On : 23 Apr 2015 19:09
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG042315

Quant Time: Apr 24 02:19:54 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 23 19:19:11 2015
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	1.571	1.605	-2.2	95	0.00
46		2-Chloronaphthalene	1.207	1.229	-1.8	95	0.00
47		2-Nitroaniline	0.329	0.358	-8.8	95	0.00
48		Acenaphthylene	2.115	2.208	-4.4	96	0.00
49		Dimethylphthalate	1.496	1.582	-5.7	98	0.00
50		2,6-Dinitrotoluene	0.367	0.392	-6.8	96	0.00
51	C	Acenaphthene	1.269	1.311	-3.3	97	0.00
52		3-Nitroaniline	0.444	0.486	-9.5	95	0.00
53	P	2,4-Dinitrophenol	0.164	0.168	-2.4	88	0.00
54		Dibenzofuran	1.790	1.836	-2.6	96	0.00
55	P	4-Nitrophenol	0.333	0.355	-6.6	96	0.00
56		2,4-Dinitrotoluene	0.503	0.550	-9.3	95	0.00
57		Fluorene	1.565	1.643	-5.0	97	0.00
58		2,3,4,6-Tetrachlorophenol	0.279	0.295	-5.7	95	0.00
59		Diethylphthalate	1.559	1.646	-5.6	96	0.00
60		4-Chlorophenyl-phenylether	0.654	0.673	-2.9	96	0.00
61		4-Nitroaniline	0.491	0.531	-8.1	93	0.00
62		Azobenzene	1.424	1.499	-5.3	96	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	98	0.00
64		4,6-Dinitro-2-methylphenol	0.135	0.140	-3.7	94	0.00
65	c	n-Nitrosodiphenylamine	0.686	0.717	-4.5	97	0.00
66		4-Bromophenyl-phenylether	0.167	0.172	-3.0	96	0.00
67		Hexachlorobenzene	0.172	0.178	-3.5	98	0.00
68		Atrazine	0.184	0.197	-7.1	98	0.00
69	C	Pentachlorophenol	0.083	0.086	-3.6	92	0.00
70		Phenanthrene	1.149	1.194	-3.9	97	0.00
71		Anthracene	1.143	1.199	-4.9	97	0.00
72		Carbazole	1.145	1.217	-6.3	98	0.00
73		Di-n-butylphthalate	1.365	1.484	-8.7	98	0.00
74	C	Fluoranthene	1.217	1.274	-4.7	98	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	98	0.00
76		Benzidine	0.681	0.691	-1.5	99	0.00
77		Pyrene	1.405	1.494	-6.3	99	0.00
78	S	Terphenyl-d14	0.868	0.914	-5.3	98	0.00
79		Butylbenzylphthalate	0.691	0.752	-8.8	97	0.00
80		Benzo(a)anthracene	1.213	1.264	-4.2	97	0.00
81		3,3'-Dichlorobenzidine	0.395	0.422	-6.8	98	0.00
82		Chrysene	1.163	1.220	-4.9	98	0.00
83		Bis(2-ethylhexyl)phthalate	0.940	1.023	-8.8	97	0.00
84	c	Di-n-octyl phthalate	1.566	1.705	-8.9	96	0.00
85		Indeno(1,2,3-cd)pyrene	1.267	1.340	-5.8	99	0.00
86	I	Perylene-d12	1.000	1.000	0.0	99	0.00
87		Benzo(b)fluoranthene	1.215	1.293	-6.4	101	0.00

Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG042315\
Data File : BG016659.D
Acq On : 23 Apr 2015 19:09
Operator : TP/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG042315

Quant Time: Apr 24 02:19:54 2015
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG042315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Apr 23 19:19:11 2015
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	1.186	1.207	-1.8	94	0.00
89 C	Benzo(a)pyrene	1.145	1.195	-4.4	98	0.00
90	Dibenzo(a,h)anthracene	1.113	1.175	-5.6	98	0.00
91	Benzo(g,h,i)perylene	1.133	1.177	-3.9	97	0.00

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