Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG120215\
Data File : BG019840.D
Acq On : 2 Dec 2015 18:41
Operator : UM/NP
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG120215

Quant Time: Dec 03 00:09:46 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 18:50:01 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	104	0.00
2		1,4-Dioxane	0.431	0.468	-8.6	114	0.00
3		Pyridine	1.309	1.454	-11.1	115	0.00
4		n-Nitrosodimethylamine	0.689	0.748	-8.6	107	0.00
5	S	2-Fluorophenol	1.107	1.183	-6.9	112	0.00
6		Aniline	2.164	2.320	-7.2	109	0.00
7	S	Phenol-d6	1.671	1.764	-5.6	110	0.00
8		2-Chlorophenol	1.349	1.416	-5.0	108	0.00
9		Benzaldehyde	1.114	1.191	-6.9	112	0.00
10	C	Phenol	1.758	1.866	-6.1	108	0.00
11		bis(2-Chloroethyl)ether	1.302	1.382	-6.1	111	0.00
12		1,3-Dichlorobenzene	1.529	1.578	-3.2	108	0.00
13	C	1,4-Dichlorobenzene	1.580	1.614	-2.2	108	0.00
14		1,2-Dichlorobenzene	1.507	1.573	-4.4	109	0.00
15		Benzyl Alcohol	1.466	1.530	-4.4	110	0.00
16		2,2'-oxybis(1-Chloropropane	2.124	2.322	-9.3	114	0.00
17		2-Methylphenol	1.241	1.302	-4.9	107	0.00
18		Hexachloroethane	0.591	0.621	-5.1	107	0.00
19	P	n-Nitroso-di-n-propylamine	1.318	1.407	-6.8	112	0.00
20		3+4-Methylphenols	1.748	1.871	-7.0	107	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
22		Acetophenone	0.548	0.576	-5.1	110	0.00
23	S	Nitrobenzene-d5	0.392	0.421	-7.4	112	0.00
24		Nitrobenzene	0.425	0.468	-10.1	113	0.00
25		Isophorone	0.840	0.890	-6.0	111	0.00
26	C	2-Nitrophenol	0.164	0.178	-8.5	111	0.00
27		2,4-Dimethylphenol	0.342	0.348	-1.8	108	0.00
28		bis(2-Chloroethoxy)methane	0.411	0.434	-5.6	112	0.00
29	C	2,4-Dichlorophenol	0.349	0.353	-1.1	104	0.00
30		1,2,4-Trichlorobenzene	0.395	0.402	-1.8	106	0.00
31		Naphthalene	1.071	1.105	-3.2	108	0.00
32		Benzoic acid	0.211	0.243	-15.2	122	0.00
33		4-Chloroaniline	0.472	0.476	-0.8	106	0.00
34	C	Hexachlorobutadiene	0.291	0.282	3.1	102	0.00
35		Caprolactam	0.130	0.134	-3.1	112	0.00
36	C	4-Chloro-3-methylphenol	0.435	0.435	0.0	106	0.00
37		2-Methylnaphthalene	0.785	0.793	-1.0	107	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	100	0.00
39		1,2,4,5-Tetrachlorobenzene	0.759	0.779	-2.6	103	0.00
40	P	Hexachlorocyclopentadiene	0.433	0.472	-9.0	106	0.00
41	S	2,4,6-Tribromophenol	0.306	0.298	2.6	98	0.00
42	C	2,4,6-Trichlorophenol	0.502	0.503	-0.2	102	0.00
43		2,4,5-Trichlorophenol	0.531	0.535	-0.8	103	0.00
44	S	2-Fluorobiphenyl	1.465	1.521	-3.8	105	0.00

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Instrument : BNA_G
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Quant Time: Dec 03 00:09:46 2015
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG120215.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 02 18:50:01 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	1.641	1.704	-3.8	104	0.00
46		2-Chloronaphthalene	1.265	1.325	-4.7	105	0.00
47		2-Nitroaniline	0.400	0.442	-10.5	109	0.00
48		Acenaphthylene	2.155	2.246	-4.2	105	0.00
49		Dimethylphthalate	1.799	1.822	-1.3	103	0.00
50		2,6-Dinitrotoluene	0.338	0.378	-11.8	107	0.00
51	C	Acenaphthene	1.308	1.366	-4.4	105	0.00
52		3-Nitroaniline	0.351	0.389	-10.8	109	0.00
53	P	2,4-Dinitrophenol	0.143	0.167	-16.8	119	0.00
54		Dibenzofuran	1.945	2.005	-3.1	104	0.00
55	P	4-Nitrophenol	0.306	0.329	-7.5	105	0.00
56		2,4-Dinitrotoluene	0.472	0.528	-11.9	106	0.00
57		Fluorene	1.741	1.771	-1.7	104	0.00
58		2,3,4,6-Tetrachlorophenol	0.491	0.481	2.0	96	0.00
59		Diethylphthalate	1.947	1.965	-0.9	104	0.00
60		4-Chlorophenyl-phenylether	1.000	1.010	-1.0	102	0.00
61		4-Nitroaniline	0.395	0.424	-7.3	105	0.00
62		Azobenzene	1.620	1.696	-4.7	106	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	97	0.00
64		4,6-Dinitro-2-methylphenol	0.103	0.114	-10.7	107	0.00
65	c	n-Nitrosodiphenylamine	0.579	0.599	-3.5	102	0.00
66		4-Bromophenyl-phenylether	0.249	0.256	-2.8	100	0.00
67		Hexachlorobenzene	0.259	0.259	0.0	98	0.00
68		Atrazine	0.251	0.255	-1.6	98	0.00
69	C	Pentachlorophenol	0.151	0.160	-6.0	102	0.00
70		Phenanthrene	1.132	1.165	-2.9	101	0.00
71		Anthracene	1.153	1.188	-3.0	101	0.00
72		Carbazole	1.015	1.062	-4.6	102	0.00
73		Di-n-butylphthalate	1.245	1.306	-4.9	102	0.00
74	C	Fluoranthene	1.447	1.497	-3.5	100	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
76		Benzidine	0.657	0.699	-6.4	105	0.00
77		Pyrene	1.177	1.178	-0.1	100	0.00
78	S	Terphenyl-d14	0.820	0.820	0.0	100	0.00
79		Butylbenzylphthalate	0.461	0.497	-7.8	108	0.00
80		Benzo(a)anthracene	1.224	1.243	-1.6	104	0.00
81		3,3'-Dichlorobenzidine	0.466	0.471	-1.1	102	0.00
82		Chrysene	1.162	1.196	-2.9	105	0.00
83		Bis(2-ethylhexyl)phthalate	0.677	0.719	-6.2	109	0.00
84	c	Di-n-octyl phthalate	1.100	1.178	-7.1	107	0.00
85		Indeno(1,2,3-cd)pyrene	1.280	1.330	-3.9	106	0.00
86	I	Perylene-d12	1.000	1.000	0.0	102	0.00
87		Benzo(b)fluoranthene	1.268	1.311	-3.4	107	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.255	1.285	-2.4	107	0.00
89 C	Benzo(a)pyrene	1.203	1.249	-3.8	107	0.00
90	Dibenzo(a,h)anthracene	1.189	1.219	-2.5	107	0.00
91	Benzo(g,h,i)perylene	1.167	1.185	-1.5	105	0.00

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