Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG022316\
Data File : BG021037.D
Acq On : 23 Feb 2016 20:29
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_G
ClientSampleId : ICVBG022316

Quant Time: Feb 24 04:05:35 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG022316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Feb 24 00:35:06 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	64	0.00
2		1,4-Dioxane	0.295	0.316	-7.1	68	0.00
3		Pyridine	1.043	1.121	-7.5	71	0.00
4		n-Nitrosodimethylamine	0.429	0.474	-10.5	69	0.00
5	S	2-Fluorophenol	1.141	1.161	-1.8	64	0.00
6		Aniline	2.072	2.256	-8.9	68	0.00
7	S	Phenol-d6	1.757	1.813	-3.2	64	0.00
8		2-Chlorophenol	1.369	1.392	-1.7	63	0.00
9		Benzaldehyde	0.850	1.055	-24.1	74	0.00
10	C	Phenol	1.833	1.925	-5.0	64	0.00
11		bis(2-Chloroethyl)ether	1.471	1.470	0.1	65	0.00
12		1,3-Dichlorobenzene	1.550	1.521	1.9	62	0.00
13	C	1,4-Dichlorobenzene	1.618	1.618	0.0	64	0.00
14		1,2-Dichlorobenzene	1.503	1.524	-1.4	65	0.00
15		Benzyl Alcohol	1.196	1.418	-18.6	73	0.00
16		2,2'-oxybis(1-Chloropropane	2.477	2.432	1.8	63	0.00
17		2-Methylphenol	1.247	1.313	-5.3	66	0.00
18		Hexachloroethane	0.581	0.579	0.3	63	0.00
19	P	n-Nitroso-di-n-propylamine	1.236	1.376	-11.3	69	0.00
20		3+4-Methylphenols	1.729	1.817	-5.1	65	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	65	0.00
22		Acetophenone	0.463	0.457	1.3	64	0.00
23	S	Nitrobenzene-d5	0.341	0.358	-5.0	65	0.00
24		Nitrobenzene	0.370	0.369	0.3	63	0.00
25		Isophorone	0.702	0.732	-4.3	65	0.00
26	C	2-Nitrophenol	0.164	0.169	-3.0	62	0.00
27		2,4-Dimethylphenol	0.296	0.303	-2.4	66	0.00
28		bis(2-Chloroethoxy)methane	0.380	0.403	-6.1	68	0.00
29	C	2,4-Dichlorophenol	0.271	0.278	-2.6	64	0.00
30		1,2,4-Trichlorobenzene	0.290	0.283	2.4	62	0.00
31		Naphthalene	1.046	1.057	-1.1	64	0.00
32		Benzoic acid	0.224	0.205	8.5	57	-0.04
33		4-Chloroaniline	0.419	0.449	-7.2	66	0.00
34	C	Hexachlorobutadiene	0.161	0.164	-1.9	63	0.00
35		Caprolactam	0.106	0.109	-2.8	63	-0.02
36	C	4-Chloro-3-methylphenol	0.343	0.363	-5.8	65	0.00
37		2-Methylnaphthalene	0.761	0.810	-6.4	67	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	64	0.00
39		1,2,4,5-Tetrachlorobenzene	0.553	0.539	2.5	61	0.00
40	P	Hexachlorocyclopentadiene	0.313	0.343	-9.6	66	0.00
41	S	2,4,6-Tribromophenol	0.240	0.242	-0.8	62	0.00
42	C	2,4,6-Trichlorophenol	0.393	0.395	-0.5	62	0.00
43		2,4,5-Trichlorophenol	0.403	0.426	-5.7	65	0.00
44	S	2-Fluorobiphenyl	1.437	1.413	1.7	62	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.735	1.675	3.5	61	0.00
46		2-Chloronaphthalene	1.214	1.191	1.9	62	0.00
47		2-Nitroaniline	0.366	0.407	-11.2	67	0.00
48		Acenaphthylene	2.219	2.186	1.5	61	0.00
49		Dimethylphthalate	1.631	1.619	0.7	62	0.00
50		2,6-Dinitrotoluene	0.334	0.356	-6.6	64	0.00
51	C	Acenaphthene	1.276	1.282	-0.5	63	0.00
52		3-Nitroaniline	0.384	0.424	-10.4	65	0.00
53	P	2,4-Dinitrophenol	0.180	0.186	-3.3	61	0.00
54		Dibenzofuran	1.812	1.990	-9.8	68	0.00
55	P	4-Nitrophenol	0.331	0.342	-3.3	65	0.00
56		2,4-Dinitrotoluene	0.480	0.477	0.6	61	0.00
57		Fluorene	1.712	1.705	0.4	62	0.00
58		2,3,4,6-Tetrachlorophenol	0.337	0.385	-14.2	69	0.00
59		Diethylphthalate	1.788	1.802	-0.8	63	0.00
60		4-Chlorophenyl-phenylether	0.789	0.765	3.0	62	0.00
61		4-Nitroaniline	0.402	0.413	-2.7	63	0.00
62		Azobenzene	1.550	1.751	-13.0	70	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	64	0.00
64		4,6-Dinitro-2-methylphenol	0.121	0.122	-0.8	60	0.00
65	c	n-Nitrosodiphenylamine	0.602	0.623	-3.5	64	0.00
66		4-Bromophenyl-phenylether	0.210	0.209	0.5	62	0.00
67		Hexachlorobenzene	0.230	0.236	-2.6	63	0.00
68		Atrazine	0.189	0.177	6.3	58	0.00
69	C	Pentachlorophenol	0.133	0.134	-0.8	61	0.00
70		Phenanthrene	1.131	1.180	-4.3	66	0.00
71		Anthracene	1.127	1.151	-2.1	64	0.00
72		Carbazole	1.016	1.063	-4.6	65	0.00
73		Di-n-butylphthalate	1.238	1.266	-2.3	63	0.00
74	C	Fluoranthene	1.138	1.168	-2.6	64	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	75	0.00
76		Benzidine	1.017	0.903	11.2	62	0.00
77		Pyrene	2.285	2.065	9.6	66	0.00
78	S	Terphenyl-d14	1.541	1.393	9.6	65	0.00
79		Butylbenzylphthalate	0.806	0.783	2.9	69	0.00
80		Benzo(a)anthracene	1.199	1.236	-3.1	76	0.00
81		3,3'-Dichlorobenzidine	0.443	0.462	-4.3	77	0.00
82		Chrysene	1.119	1.089	2.7	73	0.00
83		Bis(2-ethylhexyl)phthalate	0.880	0.887	-0.8	75	0.00
84	c	Di-n-octyl phthalate	1.252	1.292	-3.2	80	0.00
85		Indeno(1,2,3-cd)pyrene	1.006	1.036	-3.0	78	0.00
86	I	Perylene-d12	1.000	1.000	0.0	80	-0.01
87		Benzo(b)fluoranthene	1.319	1.348	-2.2	80	0.00

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88	Benzo(k)fluoranthene	1.244	1.226	1.4	77	0.00
89 C	Benzo(a)pyrene	1.181	1.226	-3.8	79	0.00
90	Dibenzo(a,h)anthracene	1.064	1.070	-0.6	78	-0.02
91	Benzo(g,h,i)perylene	1.062	1.056	0.6	78	-0.02

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