Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG122916\
Data File : BG025400.D
Acq On : 29 Dec 2016 12:12
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Dec 30 00:36:19 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 21 18:42:01 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	85	0.00
2		1,4-Dioxane	0.469	0.523	-11.5	94	0.00
3		Pyridine	1.361	1.536	-12.9	87	0.00
4		n-Nitrosodimethylamine	0.808	0.929	-15.0	90	0.00
5	S	2-Fluorophenol	1.100	1.202	-9.3	90	0.00
6		Aniline	2.100	2.314	-10.2	89	0.00
7	S	Phenol-d6	1.550	1.751	-13.0	89	0.00
8		2-Chlorophenol	1.241	1.309	-5.5	86	0.00
9		Benzaldehyde	1.134	1.161	-2.4	86	0.00
10	C	Phenol	1.718	1.979	-15.2	92	0.00
11		bis(2-Chloroethyl)ether	1.324	1.460	-10.3	90	0.00
12		1,3-Dichlorobenzene	1.502	1.638	-9.1	88	0.00
13	C	1,4-Dichlorobenzene	1.557	1.715	-10.1	90	0.00
14		1,2-Dichlorobenzene	1.486	1.601	-7.7	89	0.00
15		Benzyl Alcohol	1.479	1.712	-15.8	88	0.00
16		2,2'-oxybis(1-Chloropropane	2.451	2.702	-10.2	89	0.00
17		2-Methylphenol	1.128	1.289	-14.3	91	0.00
18		Hexachloroethane	0.598	0.657	-9.9	86	0.00
19	P	n-Nitroso-di-n-propylamine	1.309	1.482	-13.2	91	0.00
20		3+4-Methylphenols	1.561	1.835	-17.6	93	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	93	0.00
22		Acetophenone	0.556	0.565	-1.6	87	0.00
23	S	Nitrobenzene-d5	0.453	0.473	-4.4	89	0.00
24		Nitrobenzene	0.497	0.515	-3.6	89	0.00
25		Isophorone	0.820	0.854	-4.1	88	0.00
26	C	2-Nitrophenol	0.190	0.204	-7.4	91	0.00
27		2,4-Dimethylphenol	0.312	0.326	-4.5	88	0.00
28		bis(2-Chloroethoxy)methane	0.426	0.430	-0.9	91	0.00
29	C	2,4-Dichlorophenol	0.334	0.364	-9.0	93	0.00
30		1,2,4-Trichlorobenzene	0.404	0.410	-1.5	90	0.00
31		Naphthalene	0.999	1.034	-3.5	89	0.00
32		Benzoic acid	0.251	0.272	-8.4	91	-0.01
33		4-Chloroaniline	0.420	0.443	-5.5	87	0.00
34	C	Hexachlorobutadiene	0.329	0.354	-7.6	93	0.00
35		Caprolactam	0.126	0.126	0.0	87	0.00
36	C	4-Chloro-3-methylphenol	0.412	0.438	-6.3	90	0.00
37		2-Methylnaphthalene	0.740	0.793	-7.2	92	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	91	0.00
39		1,2,4,5-Tetrachlorobenzene	0.660	0.701	-6.2	91	0.00
40	P	Hexachlorocyclopentadiene	0.192	0.236	-22.9	98	0.00
41	S	2,4,6-Tribromophenol	0.322	0.336	-4.3	88	0.00
42	C	2,4,6-Trichlorophenol	0.416	0.433	-4.1	89	0.00
43		2,4,5-Trichlorophenol	0.435	0.436	-0.2	86	0.00
44	S	2-Fluorobiphenyl	1.235	1.295	-4.9	92	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.345	1.387	-3.1	89	0.00
46		2-Chloronaphthalene	1.051	1.076	-2.4	90	0.00
47		2-Nitroaniline	0.444	0.480	-8.1	90	0.00
48		Acenaphthylene	1.629	1.688	-3.6	90	0.00
49		Dimethylphthalate	1.458	1.554	-6.6	91	0.00
50		2,6-Dinitrotoluene	0.319	0.316	0.9	85	0.00
51	C	Acenaphthene	1.065	1.113	-4.5	90	0.00
52		3-Nitroaniline	0.306	0.317	-3.6	87	0.00
53	P	2,4-Dinitrophenol	0.183	0.191	-4.4	84	0.00
54		Dibenzofuran	1.684	1.769	-5.0	90	0.00
55	P	4-Nitrophenol	0.229	0.241	-5.2	86	0.00
56		2,4-Dinitrotoluene	0.464	0.482	-3.9	88	0.00
57		Fluorene	1.410	1.469	-4.2	90	0.00
58		2,3,4,6-Tetrachlorophenol	0.466	0.496	-6.4	92	0.00
59		Diethylphthalate	1.529	1.550	-1.4	87	0.00
60		4-Chlorophenyl-phenylether	0.799	0.826	-3.4	89	0.00
61		4-Nitroaniline	0.308	0.315	-2.3	79	0.00
62		Azobenzene	1.475	1.540	-4.4	89	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	89	0.00
64		4,6-Dinitro-2-methylphenol	0.139	0.150	-7.9	83	0.00
65	c	n-Nitrosodiphenylamine	0.541	0.574	-6.1	89	0.00
66		4-Bromophenyl-phenylether	0.247	0.262	-6.1	88	0.00
67		Hexachlorobenzene	0.283	0.298	-5.3	89	0.00
68		Atrazine	0.237	0.246	-3.8	87	0.00
69	C	Pentachlorophenol	0.147	0.147	0.0	80	0.00
70		Phenanthrene	1.031	1.082	-4.9	88	0.00
71		Anthracene	1.047	1.097	-4.8	89	0.00
72		Carbazole	0.937	0.974	-3.9	88	0.00
73		Di-n-butylphthalate	1.152	1.177	-2.2	86	0.00
74	C	Fluoranthene	1.361	1.416	-4.0	89	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	95	0.00
76		Benzidine	0.499	0.470	5.8	85	0.00
77		Pyrene	1.045	1.067	-2.1	89	0.00
78	S	Terphenyl-d14	0.754	0.763	-1.2	90	0.00
79		Butylbenzylphthalate	0.420	0.421	-0.2	91	0.00
80		Benzo(a)anthracene	1.116	1.126	-0.9	90	0.00
81		3,3'-Dichlorobenzidine	0.433	0.453	-4.6	92	0.00
82		Chrysene	1.070	1.104	-3.2	93	0.00
83		Bis(2-ethylhexyl)phthalate	0.584	0.591	-1.2	92	0.00
84	c	Di-n-octyl phthalate	0.970	1.005	-3.6	93	0.00
85		Indeno(1,2,3-cd)pyrene	1.360	1.432	-5.3	94	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	96	0.00
87		Benzo(b)fluoranthene	1.091	1.142	-4.7	96	0.00

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88	Benzo(k)fluoranthene	1.054	1.087	-3.1	93	0.00
89 C	Benzo(a)pyrene	1.054	1.083	-2.8	93	0.00
90	Dibenzo(a,h)anthracene	1.117	1.156	-3.5	93	-0.02
91	Benzo(g,h,i)perylene	1.110	1.151	-3.7	95	-0.03

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