Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG030217\
Data File : BG026040.D
Acq On : 2 Mar 2017 21:56
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Mar 03 00:41:48 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG021717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Feb 28 15:36:56 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	136	-0.01
2		1,4-Dioxane	0.525	0.542	-3.2	139	-0.01
3		Pyridine	1.564	1.450	7.3	121	-0.02
4		n-Nitrosodimethylamine	0.563	0.470	16.5	109	-0.01
5	S	2-Fluorophenol	1.149	1.161	-1.0	133	-0.01
6		Aniline	2.363	2.167	8.3	120	-0.01
7	S	Phenol-d6	1.700	1.571	7.6	120	0.00
8		2-Chlorophenol	1.365	1.335	2.2	129	-0.01
9		Benzaldehyde	1.007	0.882	12.4	116	-0.01
10	C	Phenol	1.847	1.702	7.9	119	0.00
11		bis(2-Chloroethyl)ether	1.518	1.398	7.9	122	-0.01
12		1,3-Dichlorobenzene	1.578	1.546	2.0	132	-0.02
13	C	1,4-Dichlorobenzene	1.599	1.565	2.1	130	-0.01
14		1,2-Dichlorobenzene	1.570	1.519	3.2	129	-0.01
15		Benzyl Alcohol	1.245	1.138	8.6	115	-0.01
16		2,2'-oxybis(1-Chloropropane	2.203	1.621	26.4#	98	-0.03
17		2-Methylphenol	1.331	1.223	8.1	118	-0.01
18		Hexachloroethane	0.530	0.540	-1.9	137	-0.02
19	P	n-Nitroso-di-n-propylamine	1.242	1.097	11.7	114	-0.01
20		3+4-Methylphenols	1.896	1.704	10.1	116	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	116	-0.01
22		Acetophenone	0.480	0.479	0.2	113	-0.01
23	S	Nitrobenzene-d5	0.317	0.331	-4.4	117	-0.01
24		Nitrobenzene	0.343	0.348	-1.5	114	-0.01
25		Isophorone	0.697	0.690	1.0	110	-0.01
26	C	2-Nitrophenol	0.160	0.178	-11.2	120	-0.01
27		2,4-Dimethylphenol	0.299	0.296	1.0	112	-0.01
28		bis(2-Chloroethoxy)methane	0.442	0.428	3.2	110	-0.02
29	C	2,4-Dichlorophenol	0.285	0.293	-2.8	114	-0.01
30		1,2,4-Trichlorobenzene	0.310	0.323	-4.2	121	-0.02
31		Naphthalene	1.034	1.027	0.7	114	-0.02
32		Benzoic acid	0.228	0.215	5.7	103	0.00
33		4-Chloroaniline	0.449	0.421	6.2	106	-0.01
34	C	Hexachlorobutadiene	0.177	0.189	-6.8	124	-0.02
35		Caprolactam	0.115	0.116	-0.9	110	-0.01
36	C	4-Chloro-3-methylphenol	0.342	0.328	4.1	107	-0.01
37		2-Methylnaphthalene	0.788	0.764	3.0	110	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	103	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.501	0.546	-9.0	110	-0.01
40	P	Hexachlorocyclopentadiene	0.274	0.292	-6.6	106	-0.02
41	S	2,4,6-Tribromophenol	0.185	0.196	-5.9	107	-0.01
42	C	2,4,6-Trichlorophenol	0.324	0.358	-10.5	109	-0.01
43		2,4,5-Trichlorophenol	0.368	0.398	-8.2	107	-0.01
44	S	2-Fluorobiphenyl	1.145	1.192	-4.1	106	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.449	1.510	-4.2	106	-0.01
46		2-Chloronaphthalene	1.048	1.087	-3.7	105	-0.01
47		2-Nitroaniline	0.332	0.335	-0.9	97	-0.01
48		Acenaphthylene	1.866	1.940	-4.0	105	-0.01
49		Dimethylphthalate	1.372	1.392	-1.5	103	-0.01
50		2,6-Dinitrotoluene	0.308	0.328	-6.5	103	-0.01
51	C	Acenaphthene	1.232	1.272	-3.2	105	-0.01
52		3-Nitroaniline	0.345	0.343	0.6	96	-0.01
53	P	2,4-Dinitrophenol	0.161	0.151	6.2	95	0.00
54		Dibenzofuran	1.643	1.642	0.1	102	-0.01
55	P	4-Nitrophenol	0.281	0.257	8.5	88	0.00
56		2,4-Dinitrotoluene	0.415	0.448	-8.0	103	-0.01
57		Fluorene	1.468	1.464	0.3	102	-0.01
58		2,3,4,6-Tetrachlorophenol	0.324	0.337	-4.0	102	0.00
59		Diethylphthalate	1.419	1.440	-1.5	103	-0.01
60		4-Chlorophenyl-phenylether	0.680	0.683	-0.4	103	-0.01
61		4-Nitroaniline	0.357	0.337	5.6	93	-0.01
62		Azobenzene	1.220	1.219	0.1	100	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	102	-0.01
64		4,6-Dinitro-2-methylphenol	0.117	0.124	-6.0	105	0.00
65	c	n-Nitrosodiphenylamine	0.607	0.611	-0.7	101	-0.01
66		4-Bromophenyl-phenylether	0.210	0.217	-3.3	103	-0.01
67		Hexachlorobenzene	0.216	0.225	-4.2	106	-0.01
68		Atrazine	0.215	0.228	-6.0	105	-0.01
69	C	Pentachlorophenol	0.133	0.142	-6.8	101	0.00
70		Phenanthrene	1.086	1.076	0.9	100	-0.01
71		Anthracene	1.107	1.110	-0.3	101	-0.01
72		Carbazole	1.112	1.075	3.3	98	-0.01
73		Di-n-butylphthalate	1.092	1.212	-11.0	108	-0.02
74	C	Fluoranthene	1.195	1.206	-0.9	102	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	99	-0.02
76		Benzidine	0.631	0.513	18.7	76	-0.01
77		Pyrene	1.265	1.321	-4.4	102	-0.01
78	S	Terphenyl-d14	0.822	0.839	-2.1	101	-0.02
79		Butylbenzylphthalate	0.480	0.584	-21.7	113	-0.02
80		Benzo(a)anthracene	1.168	1.178	-0.9	99	-0.02
81		3,3'-Dichlorobenzidine	0.416	0.421	-1.2	96	-0.02
82		Chrysene	1.123	1.140	-1.5	100	-0.02
83		Bis(2-ethylhexyl)phthalate	0.693	0.862	-24.4	116	-0.03
84	c	Di-n-octyl phthalate	1.145	1.481	-29.3#	119	-0.04
85		Indeno(1,2,3-cd)pyrene	1.323	1.181	10.7	86	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	98	-0.03
87		Benzo(b)fluoranthene	1.198	1.205	-0.6	95	-0.03

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88	Benzo(k)fluoranthene	1.169	1.194	-2.1	98	-0.03
89 C	Benzo(a)pyrene	1.134	1.113	1.9	94	-0.03
90	Dibenzo(a,h)anthracene	1.139	0.995	12.6	83	-0.06
91	Benzo(g,h,i)perylene	1.142	1.058	7.4	89	-0.06

(#) = Out of Range SPCC's out = 0 CCC's out = 1