Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071316\
Data File : BG023082.D
Acq On : 13 Jul 2016 23:51
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 14 01:04:05 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jul 12 18:30:58 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	125	0.00
2		1,4-Dioxane	0.475	0.477	-0.4	131	0.00
3		Pyridine	1.627	1.561	4.1	121	0.00
4		n-Nitrosodimethylamine	0.698	0.726	-4.0	129	0.00
5	S	2-Fluorophenol	1.223	1.208	1.2	126	0.00
6		Aniline	2.654	2.607	1.8	122	0.00
7	S	Phenol-d6	1.915	1.884	1.6	121	0.00
8		2-Chlorophenol	1.301	1.299	0.2	125	0.00
9		Benzaldehyde	1.280	1.192	6.9	119	0.00
10	C	Phenol	1.971	1.920	2.6	121	0.01
11		bis(2-Chloroethyl)ether	1.562	1.505	3.6	121	0.00
12		1,3-Dichlorobenzene	1.593	1.567	1.6	127	0.00
13	C	1,4-Dichlorobenzene	1.595	1.580	0.9	126	0.00
14		1,2-Dichlorobenzene	1.585	1.523	3.9	125	0.00
15		Benzyl Alcohol	1.754	1.695	3.4	118	0.00
16		2,2'-oxybis(1-Chloropropane	1.908	1.907	0.1	126	0.00
17		2-Methylphenol	1.283	1.214	5.4	118	0.01
18		Hexachloroethane	0.673	0.678	-0.7	136	0.00
19	P	n-Nitroso-di-n-propylamine	1.727	1.588	8.0	112	0.00
20		3+4-Methylphenols	1.789	1.743	2.6	115	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	115	0.01
22		Acetophenone	0.600	0.577	3.8	111	0.00
23	S	Nitrobenzene-d5	0.467	0.470	-0.6	117	0.00
24		Nitrobenzene	0.496	0.501	-1.0	119	0.00
25		Isophorone	0.939	0.889	5.3	106	0.00
26	C	2-Nitrophenol	0.195	0.196	-0.5	109	0.00
27		2,4-Dimethylphenol	0.337	0.320	5.0	110	0.00
28		bis(2-Chloroethoxy)methane	0.524	0.504	3.8	110	0.00
29	C	2,4-Dichlorophenol	0.359	0.348	3.1	109	0.00
30		1,2,4-Trichlorobenzene	0.411	0.401	2.4	113	0.01
31		Naphthalene	1.018	1.010	0.8	115	0.00
32		Benzoic acid	0.306	0.287	6.2	102	0.03
33		4-Chloroaniline	0.498	0.482	3.2	111	0.01
34	C	Hexachlorobutadiene	0.334	0.320	4.2	113	0.00
35		Caprolactam	0.148	0.125	15.5	97	0.02
36	C	4-Chloro-3-methylphenol	0.491	0.459	6.5	105	0.01
37		2-Methylnaphthalene	0.805	0.782	2.9	109	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	103	0.01
39		1,2,4,5-Tetrachlorobenzene	0.862	0.851	1.3	103	0.00
40	P	Hexachlorocyclopentadiene	0.496	0.595	-20.0	125	0.00
41	S	2,4,6-Tribromophenol	0.359	0.311	13.4	92	0.00
42	C	2,4,6-Trichlorophenol	0.495	0.480	3.0	99	0.01
43		2,4,5-Trichlorophenol	0.509	0.486	4.5	101	0.01
44	S	2-Fluorobiphenyl	1.270	1.235	2.8	103	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.501	1.502	-0.1	104	0.01
46		2-Chloronaphthalene	1.217	1.247	-2.5	106	0.01
47		2-Nitroaniline	0.519	0.543	-4.6	109	0.02
48		Acenaphthylene	1.826	1.811	0.8	104	0.01
49		Dimethylphthalate	1.634	1.507	7.8	98	0.01
50		2,6-Dinitrotoluene	0.367	0.356	3.0	101	0.01
51	C	Acenaphthene	1.193	1.169	2.0	103	0.01
52		3-Nitroaniline	0.366	0.369	-0.8	104	0.01
53	P	2,4-Dinitrophenol	0.252	0.360	-42.9#	143	0.01
54		Dibenzofuran	1.786	1.709	4.3	102	0.01
55	P	4-Nitrophenol	0.307	0.283	7.8	96	0.02
56		2,4-Dinitrotoluene	0.535	0.512	4.3	101	0.01
57		Fluorene	1.537	1.477	3.9	103	0.00
58		2,3,4,6-Tetrachlorophenol	0.509	0.460	9.6	94	0.01
59		Diethylphthalate	1.662	1.541	7.3	98	0.01
60		4-Chlorophenyl-phenylether	0.936	0.866	7.5	96	0.01
61		4-Nitroaniline	0.396	0.389	1.8	104	0.01
62		Azobenzene	1.590	1.618	-1.8	107	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	101	0.00
64		4,6-Dinitro-2-methylphenol	0.149	0.141	5.4	91	0.01
65	c	n-Nitrosodiphenylamine	0.576	0.580	-0.7	100	0.00
66		4-Bromophenyl-phenylether	0.260	0.237	8.8	91	0.00
67		Hexachlorobenzene	0.283	0.265	6.4	94	0.00
68		Atrazine	0.256	0.243	5.1	97	0.00
69	C	Pentachlorophenol	0.183	0.170	7.1	88	0.00
70		Phenanthrene	0.980	0.943	3.8	98	0.00
71		Anthracene	0.992	0.971	2.1	100	0.00
72		Carbazole	0.858	0.837	2.4	100	0.00
73		Di-n-butylphthalate	0.954	0.993	-4.1	103	0.00
74	C	Fluoranthene	1.203	1.107	8.0	98	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	97	0.02
76		Benzidine	0.573	0.774	-35.1#	127	0.01
77		Pyrene	1.124	1.087	3.3	99	0.00
78	S	Terphenyl-d14	0.646	0.646	0.0	97	0.00
79		Butylbenzylphthalate	0.445	0.470	-5.6	103	0.01
80		Benzo(a)anthracene	1.119	1.100	1.7	98	0.02
81		3,3'-Dichlorobenzidine	0.491	0.478	2.6	93	0.02
82		Chrysene	1.050	1.026	2.3	98	0.02
83		Bis(2-ethylhexyl)phthalate	0.618	0.637	-3.1	102	0.02
84	c	Di-n-octyl phthalate	1.031	1.075	-4.3	102	0.04
85		Indeno(1,2,3-cd)pyrene	1.338	1.199	10.4	90	0.12
86	I	Perylene-d12	1.000	1.000	0.0	94	0.06
87		Benzo(b)fluoranthene	1.154	1.145	0.8	96	0.05

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88	Benzo(k)fluoranthene	1.095	1.091	0.4	92	0.06
89 C	Benzo(a)pyrene	1.106	1.094	1.1	94	0.06
90	Dibenzo(a,h)anthracene	1.098	1.023	6.8	89	0.12
91	Benzo(g,h,i)perylene	1.099	1.024	6.8	88	0.13

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