Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG073116\
Data File : BG023366.D
Acq On : 31 Jul 2016 12:52
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 01 02:13:19 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG072616.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 27 17:30:33 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	119	0.00
2		1,4-Dioxane	0.502	0.463	7.8	109	0.00
3		Pyridine	1.759	1.615	8.2	87	-0.03
4		n-Nitrosodimethylamine	0.795	0.578	27.3#	68	-0.01
5	S	2-Fluorophenol	1.172	1.185	-1.1	118	0.00
6		Aniline	2.402	2.600	-8.2	128	0.00
7	S	Phenol-d6	1.799	1.792	0.4	113	0.00
8		2-Chlorophenol	1.307	1.387	-6.1	125	0.00
9		Benzaldehyde	1.292	1.379	-6.7	125	0.00
10	C	Phenol	1.898	1.918	-1.1	117	0.00
11		bis(2-Chloroethyl)ether	1.613	1.623	-0.6	119	0.00
12		1,3-Dichlorobenzene	1.473	1.535	-4.2	125	0.00
13	C	1,4-Dichlorobenzene	1.503	1.602	-6.6	125	0.00
14		1,2-Dichlorobenzene	1.450	1.543	-6.4	126	0.00
15		Benzyl Alcohol	1.210	1.358	-12.2	127	0.00
16		2,2'-oxybis(1-Chloropropane	2.567	2.286	10.9	117	-0.02
17		2-Methylphenol	1.297	1.323	-2.0	123	0.00
18		Hexachloroethane	0.626	0.618	1.3	116	-0.02
19	P	n-Nitroso-di-n-propylamine	1.618	1.598	1.2	113	0.00
20		3+4-Methylphenols	1.764	1.869	-6.0	123	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	122	0.00
22		Acetophenone	0.522	0.518	0.8	117	0.00
23	S	Nitrobenzene-d5	0.409	0.379	7.3	111	0.00
24		Nitrobenzene	0.475	0.409	13.9	105	0.00
25		Isophorone	0.845	0.792	6.3	111	0.00
26	C	2-Nitrophenol	0.179	0.192	-7.3	127	-0.02
27		2,4-Dimethylphenol	0.302	0.311	-3.0	126	-0.02
28		bis(2-Chloroethoxy)methane	0.474	0.483	-1.9	126	0.00
29	C	2,4-Dichlorophenol	0.293	0.317	-8.2	126	-0.02
30		1,2,4-Trichlorobenzene	0.323	0.339	-5.0	124	0.00
31		Naphthalene	0.929	0.941	-1.3	124	-0.02
32		Benzoic acid	0.215	0.276	-28.4#	160#	0.00
33		4-Chloroaniline	0.429	0.480	-11.9	140	0.00
34	C	Hexachlorobutadiene	0.212	0.215	-1.4	111	0.00
35		Caprolactam	0.124	0.129	-4.0	130	0.02
36	C	4-Chloro-3-methylphenol	0.385	0.391	-1.6	118	0.00
37		2-Methylnaphthalene	0.689	0.731	-6.1	127	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	115	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.648	0.703	-8.5	122	-0.02
40	P	Hexachlorocyclopentadiene	0.307	0.367	-19.5	120	-0.02
41	S	2,4,6-Tribromophenol	0.207	0.262	-26.6#	130	-0.02
42	C	2,4,6-Trichlorophenol	0.380	0.433	-13.9	124	-0.02
43		2,4,5-Trichlorophenol	0.403	0.440	-9.2	120	-0.02
44	S	2-Fluorobiphenyl	1.069	1.111	-3.9	119	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.395	1.487	-6.6	125	0.00
46		2-Chloronaphthalene	1.118	1.162	-3.9	120	-0.01
47		2-Nitroaniline	0.473	0.458	3.2	112	-0.01
48		Acenaphthylene	1.711	1.769	-3.4	119	-0.02
49		Dimethylphthalate	1.476	1.442	2.3	112	0.00
50		2,6-Dinitrotoluene	0.346	0.352	-1.7	118	0.00
51	C	Acenaphthene	1.110	1.119	-0.8	118	-0.02
52		3-Nitroaniline	0.348	0.390	-12.1	134	-0.01
53	P	2,4-Dinitrophenol	0.202	0.231	-14.4	127	-0.01
54		Dibenzofuran	1.571	1.598	-1.7	117	-0.02
55	P	4-Nitrophenol	0.302	0.307	-1.7	123	-0.03
56		2,4-Dinitrotoluene	0.487	0.487	0.0	115	-0.01
57		Fluorene	1.387	1.388	-0.1	115	-0.02
58		2,3,4,6-Tetrachlorophenol	0.342	0.378	-10.5	119	-0.02
59		Diethylphthalate	1.512	1.455	3.8	110	-0.02
60		4-Chlorophenyl-phenylether	0.751	0.755	-0.5	114	-0.01
61		4-Nitroaniline	0.365	0.394	-7.9	125	-0.01
62		Azobenzene	1.567	1.409	10.1	104	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	113	-0.02
64		4,6-Dinitro-2-methylphenol	0.128	0.145	-13.3	121	-0.01
65	c	n-Nitrosodiphenylamine	0.534	0.578	-8.2	120	-0.01
66		4-Bromophenyl-phenylether	0.202	0.232	-14.9	124	-0.02
67		Hexachlorobenzene	0.203	0.247	-21.7	133	-0.02
68		Atrazine	0.205	0.201	2.0	110	-0.02
69	C	Pentachlorophenol	0.125	0.146	-16.8	123	-0.02
70		Phenanthrene	0.875	0.912	-4.2	113	-0.02
71		Anthracene	0.884	0.927	-4.9	112	-0.02
72		Carbazole	0.810	0.821	-1.4	116	-0.02
73		Di-n-butylphthalate	0.903	0.901	0.2	109	-0.02
74	C	Fluoranthene	0.998	0.928	7.0	110	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	105	-0.02
76		Benzidine	0.563	0.513	8.9	95	-0.02
77		Pyrene	1.135	1.165	-2.6	113	-0.02
78	S	Terphenyl-d14	0.615	0.644	-4.7	115	-0.02
79		Butylbenzylphthalate	0.499	0.495	0.8	107	-0.02
80		Benzo(a)anthracene	1.047	1.061	-1.3	109	-0.02
81		3,3'-Dichlorobenzidine	0.425	0.455	-7.1	111	-0.03
82		Chrysene	1.012	0.991	2.1	106	-0.03
83		Bis(2-ethylhexyl)phthalate	0.699	0.693	0.9	106	-0.03
84	c	Di-n-octyl phthalate	1.143	1.178	-3.1	108	-0.04
85		Indeno(1,2,3-cd)pyrene	1.180	1.297	-9.9	114	-0.08
86	I	Perylene-d12	1.000	1.000	0.0	109	-0.05
87		Benzo(b)fluoranthene	1.062	1.105	-4.0	113	-0.04

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.062	1.087	-2.4	112	-0.04
89 C	Benzo(a)pyrene	1.041	1.063	-2.1	111	-0.04
90	Dibenzo(a,h)anthracene	0.985	1.085	-10.2	119	-0.07
91	Benzo(g,h,i)perylene	1.037	1.093	-5.4	100	-0.09

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