Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG070816\
Data File : BG022916.D
Acq On : 8 Jul 2016 4:43
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 08 05:54:10 2016
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG070816.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jul 07 16:51:07 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	105	0.00
2		1,4-Dioxane	0.475	0.424	10.7	98	0.00
3		Pyridine	1.627	1.562	4.0	101	0.00
4		n-Nitrosodimethylamine	0.698	0.681	2.4	102	0.00
5	S	2-Fluorophenol	1.223	1.204	1.6	106	0.00
6		Aniline	2.654	2.644	0.4	104	0.00
7	S	Phenol-d6	1.915	1.914	0.1	103	0.00
8		2-Chlorophenol	1.301	1.299	0.2	105	0.00
9		Benzaldehyde	1.280	1.258	1.7	105	0.00
10	C	Phenol	1.971	1.982	-0.6	105	0.00
11		bis(2-Chloroethyl)ether	1.562	1.519	2.8	103	0.00
12		1,3-Dichlorobenzene	1.593	1.576	1.1	107	0.00
13	C	1,4-Dichlorobenzene	1.595	1.601	-0.4	107	0.00
14		1,2-Dichlorobenzene	1.585	1.507	4.9	104	0.00
15		Benzyl Alcohol	1.754	1.745	0.5	102	0.00
16		2,2'-oxybis(1-Chloropropane	1.908	1.839	3.6	102	0.00
17		2-Methylphenol	1.283	1.298	-1.2	106	0.00
18		Hexachloroethane	0.673	0.640	4.9	108	0.00
19	P	n-Nitroso-di-n-propylamine	1.727	1.662	3.8	99	0.00
20		3+4-Methylphenols	1.789	1.862	-4.1	103	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	102	0.00
22		Acetophenone	0.600	0.584	2.7	99	0.00
23	S	Nitrobenzene-d5	0.467	0.460	1.5	101	0.00
24		Nitrobenzene	0.496	0.490	1.2	103	0.00
25		Isophorone	0.939	0.898	4.4	95	0.00
26	C	2-Nitrophenol	0.195	0.197	-1.0	97	0.00
27		2,4-Dimethylphenol	0.337	0.327	3.0	100	0.00
28		bis(2-Chloroethoxy)methane	0.524	0.506	3.4	97	0.00
29	C	2,4-Dichlorophenol	0.359	0.363	-1.1	101	0.00
30		1,2,4-Trichlorobenzene	0.411	0.405	1.5	101	0.00
31		Naphthalene	1.018	1.003	1.5	101	0.00
32		Benzoic acid	0.306	0.291	4.9	91	0.01
33		4-Chloroaniline	0.498	0.484	2.8	99	0.00
34	C	Hexachlorobutadiene	0.334	0.326	2.4	102	0.00
35		Caprolactam	0.148	0.132	10.8	90	0.00
36	C	4-Chloro-3-methylphenol	0.491	0.472	3.9	96	0.00
37		2-Methylnaphthalene	0.805	0.778	3.4	96	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
39		1,2,4,5-Tetrachlorobenzene	0.862	0.897	-4.1	98	0.00
40	P	Hexachlorocyclopentadiene	0.496	0.510	-2.8	96	0.00
41	S	2,4,6-Tribromophenol	0.359	0.335	6.7	89	0.00
42	C	2,4,6-Trichlorophenol	0.495	0.499	-0.8	93	0.00
43		2,4,5-Trichlorophenol	0.509	0.507	0.4	95	0.00
44	S	2-Fluorobiphenyl	1.270	1.280	-0.8	97	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.501	1.533	-2.1	96	0.00
46		2-Chloronaphthalene	1.217	1.221	-0.3	94	0.00
47		2-Nitroaniline	0.519	0.508	2.1	92	0.00
48		Acenaphthylene	1.826	1.819	0.4	94	0.00
49		Dimethylphthalate	1.634	1.558	4.7	91	0.00
50		2,6-Dinitrotoluene	0.367	0.358	2.5	92	0.00
51	C	Acenaphthene	1.193	1.355	-13.6	108	0.00
52		3-Nitroaniline	0.366	0.352	3.8	90	0.00
53	P	2,4-Dinitrophenol	0.252	0.223	11.5	80	0.00
54		Dibenzofuran	1.786	1.746	2.2	94	0.00
55	P	4-Nitrophenol	0.307	0.288	6.2	88	0.00
56		2,4-Dinitrotoluene	0.535	0.500	6.5	89	0.00
57		Fluorene	1.537	1.492	2.9	93	0.00
58		2,3,4,6-Tetrachlorophenol	0.509	0.487	4.3	90	0.00
59		Diethylphthalate	1.662	1.553	6.6	89	0.00
60		4-Chlorophenyl-phenylether	0.936	0.902	3.6	90	0.00
61		4-Nitroaniline	0.396	0.373	5.8	90	0.00
62		Azobenzene	1.590	1.541	3.1	92	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
64		4,6-Dinitro-2-methylphenol	0.149	0.146	2.0	84	0.00
65	c	n-Nitrosodiphenylamine	0.576	0.586	-1.7	90	0.00
66		4-Bromophenyl-phenylether	0.260	0.260	0.0	89	0.00
67		Hexachlorobenzene	0.283	0.287	-1.4	90	0.00
68		Atrazine	0.256	0.250	2.3	89	0.00
69	C	Pentachlorophenol	0.183	0.176	3.8	82	0.00
70		Phenanthrene	0.980	0.976	0.4	90	0.00
71		Anthracene	0.992	0.992	0.0	91	0.00
72		Carbazole	0.858	0.850	0.9	90	0.00
73		Di-n-butylphthalate	0.954	0.998	-4.6	92	0.00
74	C	Fluoranthene	1.203	1.144	4.9	90	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	88	0.00
76		Benzidine	0.573	0.495	13.6	73	0.00
77		Pyrene	1.124	1.104	1.8	91	0.00
78	S	Terphenyl-d14	0.646	0.685	-6.0	93	0.00
79		Butylbenzylphthalate	0.445	0.450	-1.1	89	0.00
80		Benzo(a)anthracene	1.119	1.129	-0.9	91	0.00
81		3,3'-Dichlorobenzidine	0.491	0.492	-0.2	87	0.00
82		Chrysene	1.050	1.073	-2.2	92	0.00
83		Bis(2-ethylhexyl)phthalate	0.618	0.618	0.0	90	0.00
84	c	Di-n-octyl phthalate	1.031	1.036	-0.5	89	0.00
85		Indeno(1,2,3-cd)pyrene	1.338	1.304	2.5	88	0.01
86	I	Perylene-d12	1.000	1.000	0.0	88	0.00
87		Benzo(b)fluoranthene	1.154	1.145	0.8	90	0.00

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88	Benzo(k)fluoranthene	1.095	1.106	-1.0	88	0.00
89 C	Benzo(a)pyrene	1.106	1.098	0.7	89	0.00
90	Dibenzo(a,h)anthracene	1.098	1.088	0.9	89	0.00
91	Benzo(g,h,i)perylene	1.099	1.095	0.4	89	0.01

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