Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG071116\
Data File : BG023004.D
Acq On : 11 Jul 2016 17:18
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 12 03:49:31 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	77	0.00
2		1,4-Dioxane	0.475	0.478	-0.6	80	0.00
3		Pyridine	1.627	1.646	-1.2	78	0.00
4		n-Nitrosodimethylamine	0.698	0.702	-0.6	76	0.00
5	S	2-Fluorophenol	1.223	1.206	1.4	77	0.00
6		Aniline	2.654	2.812	-6.0	80	0.00
7	S	Phenol-d6	1.915	2.070	-8.1	81	0.00
8		2-Chlorophenol	1.301	1.384	-6.4	82	0.00
9		Benzaldehyde	1.280	1.304	-1.9	80	0.00
10	C	Phenol	1.971	2.150	-9.1	83	0.00
11		bis(2-Chloroethyl)ether	1.562	1.577	-1.0	78	0.00
12		1,3-Dichlorobenzene	1.593	1.627	-2.1	81	0.00
13	C	1,4-Dichlorobenzene	1.595	1.686	-5.7	82	0.00
14		1,2-Dichlorobenzene	1.585	1.617	-2.0	81	0.00
15		Benzyl Alcohol	1.754	1.880	-7.2	80	0.00
16		2,2'-oxybis(1-Chloropropane	1.908	2.012	-5.5	81	0.00
17		2-Methylphenol	1.283	1.381	-7.6	82	0.01
18		Hexachloroethane	0.673	0.676	-0.4	83	0.00
19	P	n-Nitroso-di-n-propylamine	1.727	1.768	-2.4	76	0.00
20		3+4-Methylphenols	1.789	1.984	-10.9	80	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	80	0.00
22		Acetophenone	0.600	0.577	3.8	77	0.00
23	S	Nitrobenzene-d5	0.467	0.473	-1.3	82	0.00
24		Nitrobenzene	0.496	0.503	-1.4	83	0.00
25		Isophorone	0.939	0.904	3.7	75	0.00
26	C	2-Nitrophenol	0.195	0.191	2.1	74	0.00
27		2,4-Dimethylphenol	0.337	0.330	2.1	79	0.00
28		bis(2-Chloroethoxy)methane	0.524	0.511	2.5	77	0.00
29	C	2,4-Dichlorophenol	0.359	0.373	-3.9	81	0.00
30		1,2,4-Trichlorobenzene	0.411	0.406	1.2	80	0.00
31		Naphthalene	1.018	1.042	-2.4	83	0.00
32		Benzoic acid	0.306	0.280	8.5	69	0.00
33		4-Chloroaniline	0.498	0.490	1.6	79	0.00
34	C	Hexachlorobutadiene	0.334	0.318	4.8	78	0.00
35		Caprolactam	0.148	0.129	12.8	69	0.00
36	C	4-Chloro-3-methylphenol	0.491	0.495	-0.8	79	0.00
37		2-Methylnaphthalene	0.805	0.823	-2.2	80	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	77	0.00
39		1,2,4,5-Tetrachlorobenzene	0.862	0.888	-3.0	80	0.00
40	P	Hexachlorocyclopentadiene	0.496	0.538	-8.5	84	0.00
41	S	2,4,6-Tribromophenol	0.359	0.294	18.1	65	0.00
42	C	2,4,6-Trichlorophenol	0.495	0.489	1.2	75	0.00
43		2,4,5-Trichlorophenol	0.509	0.492	3.3	76	0.00
44	S	2-Fluorobiphenyl	1.270	1.300	-2.4	81	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.501	1.567	-4.4	81	0.00
46		2-Chloronaphthalene	1.217	1.268	-4.2	81	0.00
47		2-Nitroaniline	0.519	0.525	-1.2	79	0.00
48		Acenaphthylene	1.826	1.862	-2.0	79	0.00
49		Dimethylphthalate	1.634	1.531	6.3	74	0.00
50		2,6-Dinitrotoluene	0.367	0.342	6.8	73	0.00
51	C	Acenaphthene	1.193	1.191	0.2	78	0.00
52		3-Nitroaniline	0.366	0.350	4.4	74	0.00
53	P	2,4-Dinitrophenol	0.252	0.265	-5.2	79	0.00
54		Dibenzofuran	1.786	1.721	3.6	76	0.00
55	P	4-Nitrophenol	0.307	0.264	14.0	67	0.00
56		2,4-Dinitrotoluene	0.535	0.492	8.0	72	0.00
57		Fluorene	1.537	1.462	4.9	76	0.00
58		2,3,4,6-Tetrachlorophenol	0.509	0.442	13.2	67	0.00
59		Diethylphthalate	1.662	1.527	8.1	73	0.00
60		4-Chlorophenyl-phenylether	0.936	0.876	6.4	72	0.00
61		4-Nitroaniline	0.396	0.345	12.9	69	0.00
62		Azobenzene	1.590	1.591	-0.1	79	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	70	0.00
64		4,6-Dinitro-2-methylphenol	0.149	0.123	17.4	55	0.00
65	c	n-Nitrosodiphenylamine	0.576	0.609	-5.7	73	0.00
66		4-Bromophenyl-phenylether	0.260	0.257	1.2	69	0.00
67		Hexachlorobenzene	0.283	0.273	3.5	67	0.00
68		Atrazine	0.256	0.248	3.1	69	0.00
69	C	Pentachlorophenol	0.183	0.155	15.3	56	0.00
70		Phenanthrene	0.980	0.986	-0.6	71	0.00
71		Anthracene	0.992	1.013	-2.1	73	0.00
72		Carbazole	0.858	0.859	-0.1	71	0.00
73		Di-n-butylphthalate	0.954	1.026	-7.5	74	0.00
74	C	Fluoranthene	1.203	1.183	1.7	73	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	69	0.01
76		Benzidine	0.573	0.862	-50.4#	101	0.00
77		Pyrene	1.124	1.132	-0.7	74	0.00
78	S	Terphenyl-d14	0.646	0.709	-9.8	76	0.00
79		Butylbenzylphthalate	0.445	0.459	-3.1	72	0.00
80		Benzo(a)anthracene	1.119	1.147	-2.5	72	0.01
81		3,3'-Dichlorobenzidine	0.491	0.489	0.4	68	0.01
82		Chrysene	1.050	1.085	-3.3	73	0.01
83		Bis(2-ethylhexyl)phthalate	0.618	0.639	-3.4	73	0.02
84	c	Di-n-octyl phthalate	1.031	1.062	-3.0	72	0.03
85		Indeno(1,2,3-cd)pyrene	1.338	1.245	7.0	66	0.09
86	I	Perylene-d12	1.000	1.000	0.0	67	0.06
87		Benzo(b)fluoranthene	1.154	1.166	-1.0	70	0.04

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88	Benzo(k)fluoranthene	1.095	1.144	-4.5	70	0.04
89 C	Benzo(a)pyrene	1.106	1.115	-0.8	69	0.05
90	Dibenzo(a,h)anthracene	1.098	1.071	2.5	67	0.09
91	Benzo(g,h,i)perylene	1.099	1.056	3.9	65	0.10

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