Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\DATA\BG070317\
Data File : BG027830.D
Acq On : 3 Jul 2017 13:36
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jul 04 06:52:26 2017
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG062817.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 28 19:26:37 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	85	-0.03
2		1,4-Dioxane	0.484	0.418	13.6	73	-0.03
3		Pyridine	1.489	1.292	13.2	71	-0.03
4		n-Nitrosodimethylamine	0.730	0.706	3.3	84	-0.03
5	S	2-Fluorophenol	1.299	1.259	3.1	80	-0.03
6		Aniline	2.300	1.769	23.1	62	-0.02
7	S	Phenol-d6	1.984	1.886	4.9	78	-0.02
8		2-Chlorophenol	1.451	1.443	0.6	83	-0.03
9		Benzaldehyde	1.181	1.039	12.0	72	-0.03
10	C	Phenol	1.963	1.819	7.3	75	-0.02
11		bis(2-Chloroethyl)ether	1.491	1.676	-12.4	93	-0.03
12		1,3-Dichlorobenzene	1.544	1.603	-3.8	86	-0.02
13	C	1,4-Dichlorobenzene	1.600	1.679	-4.9	88	-0.03
14		1,2-Dichlorobenzene	1.551	1.662	-7.2	89	-0.03
15		Benzyl Alcohol	1.373	0.582	57.6#	34#	-0.02
16		2,2'-oxybis(1-Chloropropane	3.180	2.834	10.9	74	-0.02
17		2-Methylphenol	1.389	1.391	-0.1	80	-0.02
18		Hexachloroethane	0.563	0.569	-1.1	86	-0.03
19	P	n-Nitroso-di-n-propylamine	1.443	1.317	8.7	74	-0.02
20		3+4-Methylphenols	2.003	1.873	6.5	75	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	81	-0.02
22		Acetophenone	0.483	0.484	-0.2	81	-0.03
23	S	Nitrobenzene-d5	0.359	0.363	-1.1	81	-0.02
24		Nitrobenzene	0.375	0.379	-1.1	81	-0.02
25		Isophorone	0.782	0.730	6.6	75	-0.02
26	C	2-Nitrophenol	0.169	0.179	-5.9	82	-0.03
27		2,4-Dimethylphenol	0.317	0.314	0.9	78	-0.02
28		bis(2-Chloroethoxy)methane	0.443	0.401	9.5	73	-0.02
29	C	2,4-Dichlorophenol	0.278	0.299	-7.6	85	-0.01
30		1,2,4-Trichlorobenzene	0.284	0.315	-10.9	90	-0.03
31		Naphthalene	1.064	1.078	-1.3	81	-0.02
32		Benzoic acid	0.169	0.102	39.6#	48#	-0.01
33		4-Chloroaniline	0.459	0.428	6.8	73	0.04
34	C	Hexachlorobutadiene	0.158	0.203	-28.5#	106	-0.02
35		Caprolactam	0.140	0.116	17.1	66	0.00
36	C	4-Chloro-3-methylphenol	0.391	0.363	7.2	74	-0.01
37		2-Methylnaphthalene	0.792	0.799	-0.9	81	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	77	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.526	0.639	-21.5	93	-0.02
40	P	Hexachlorocyclopentadiene	0.248	0.311	-25.4#	96	-0.02
41	S	2,4,6-Tribromophenol	0.274	0.361	-31.8#	99	-0.02
42	C	2,4,6-Trichlorophenol	0.380	0.432	-13.7	86	-0.02
43		2,4,5-Trichlorophenol	0.430	0.484	-12.6	83	-0.01
44	S	2-Fluorobiphenyl	1.400	1.593	-13.8	87	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.603	1.712	-6.8	82	-0.03
46		2-Chloronaphthalene	1.211	1.292	-6.7	82	-0.02
47		2-Nitroaniline	0.485	0.467	3.7	71	-0.02
48		Acenaphthylene	2.213	2.309	-4.3	79	-0.02
49		Dimethylphthalate	1.706	1.740	-2.0	78	-0.02
50		2,6-Dinitrotoluene	0.371	0.386	-4.0	77	-0.02
51	C	Acenaphthene	1.418	1.498	-5.6	80	-0.02
52		3-Nitroaniline	0.436	0.410	6.0	70	0.00
53	P	2,4-Dinitrophenol	0.164	0.197	-20.1	91	-0.02
54		Dibenzofuran	1.875	1.960	-4.5	80	-0.02
55	P	4-Nitrophenol	0.333	0.291	12.6	64	0.00
56		2,4-Dinitrotoluene	0.517	0.537	-3.9	77	-0.01
57		Fluorene	1.818	1.894	-4.2	79	-0.02
58		2,3,4,6-Tetrachlorophenol	0.371	0.402	-8.4	83	-0.02
59		Diethylphthalate	1.889	1.871	1.0	76	-0.02
60		4-Chlorophenyl-phenylether	0.805	0.886	-10.1	83	-0.02
61		4-Nitroaniline	0.455	0.429	5.7	71	-0.01
62		Azobenzene	1.660	1.504	9.4	69	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	77	-0.02
64		4,6-Dinitro-2-methylphenol	0.115	0.130	-13.0	86	-0.02
65	c	n-Nitrosodiphenylamine	0.651	0.674	-3.5	78	-0.02
66		4-Bromophenyl-phenylether	0.209	0.250	-19.6	90	-0.02
67		Hexachlorobenzene	0.222	0.267	-20.3	92	-0.02
68		Atrazine	0.205	0.218	-6.3	79	-0.01
69	C	Pentachlorophenol	0.123	0.143	-16.3	88	-0.02
70		Phenanthrene	1.155	1.206	-4.4	79	-0.02
71		Anthracene	1.166	1.229	-5.4	79	-0.02
72		Carbazole	1.148	1.187	-3.4	78	-0.02
73		Di-n-butylphthalate	1.262	1.299	-2.9	77	-0.02
74	C	Fluoranthene	1.262	1.358	-7.6	82	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	82	-0.03
76		Benzidine	0.447	0.258	42.3#	43#	0.00
77		Pyrene	1.255	1.288	-2.6	83	-0.02
78	S	Terphenyl-d14	0.853	0.944	-10.7	87	-0.01
79		Butylbenzylphthalate	0.565	0.534	5.5	76	-0.02
80		Benzo(a)anthracene	1.209	1.245	-3.0	83	-0.03
81		3,3'-Dichlorobenzidine	0.439	0.463	-5.5	84	-0.03
82		Chrysene	1.133	1.173	-3.5	83	-0.03
83		Bis(2-ethylhexyl)phthalate	0.826	0.774	6.3	75	-0.03
84	c	Di-n-octyl phthalate	1.365	1.297	5.0	75	-0.03
85		Indeno(1,2,3-cd)pyrene	1.295	1.453	-12.2	88	-0.07
86	I	Perylene-d12	1.000	1.000	0.0	89	-0.04
87		Benzo(b)fluoranthene	1.289	1.281	0.6	87	-0.04

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Max. RRF Dev : 25% Max. Rel. Area : 150%

	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.274	1.252	1.7	85	-0.04
89 C	Benzo(a)pyrene	1.209	1.203	0.5	86	-0.04
90	Dibenzo(a,h)anthracene	1.225	1.267	-3.4	90	-0.07
91	Benzo(g,h,i)perylene	1.177	1.218	-3.5	90	-0.07

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