Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA G\DATA\BG010917\
Data File : BG025433.D
Acq On : 9 Jan 2017 16:58
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Jan 10 00:11:17 2017
Quant Method : Z:\HPCHEM1\BNA G\METHODS\8270-BG122116.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 21 18:42:01 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	96	-0.01
2		1,4-Dioxane	0.469	0.481	-2.6	98	-0.02
3		Pyridine	1.361	1.487	-9.3	96	-0.01
4		n-Nitrosodimethylamine	0.808	0.888	-9.9	97	-0.01
5	S	2-Fluorophenol	1.100	1.197	-8.8	102	-0.02
6		Aniline	2.100	2.290	-9.0	100	0.00
7	S	Phenol-d6	1.550	1.693	-9.2	97	-0.02
8		2-Chlorophenol	1.241	1.317	-6.1	97	-0.02
9		Benzaldehyde	1.134	1.109	2.2	93	-0.02
10	C	Phenol	1.718	1.848	-7.6	97	-0.01
11		bis(2-Chloroethyl)ether	1.324	1.396	-5.4	98	-0.02
12		1,3-Dichlorobenzene	1.502	1.572	-4.7	96	0.00
13	C	1,4-Dichlorobenzene	1.557	1.617	-3.9	96	-0.01
14		1,2-Dichlorobenzene	1.486	1.509	-1.5	95	-0.02
15		Benzyl Alcohol	1.479	1.581	-6.9	93	0.00
16		2,2'-oxybis(1-Chloropropane	2.451	2.676	-9.2	100	-0.01
17		2-Methylphenol	1.128	1.241	-10.0	100	-0.02
18		Hexachloroethane	0.598	0.654	-9.4	96	-0.01
19	P	n-Nitroso-di-n-propylamine	1.309	1.401	-7.0	98	-0.01
20		3+4-Methylphenols	1.561	1.714	-9.8	99	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	100	-0.01
22		Acetophenone	0.556	0.564	-1.4	94	-0.01
23	S	Nitrobenzene-d5	0.453	0.471	-4.0	96	-0.01
24		Nitrobenzene	0.497	0.520	-4.6	98	-0.02
25		Isophorone	0.820	0.865	-5.5	97	-0.01
26	C	2-Nitrophenol	0.190	0.203	-6.8	98	-0.01
27		2,4-Dimethylphenol	0.312	0.327	-4.8	96	-0.02
28		bis(2-Chloroethoxy)methane	0.426	0.439	-3.1	101	-0.01
29	C	2,4-Dichlorophenol	0.334	0.366	-9.6	101	-0.01
30		1,2,4-Trichlorobenzene	0.404	0.409	-1.2	97	-0.01
31		Naphthalene	0.999	1.008	-0.9	94	-0.01
32		Benzoic acid	0.251	0.227	9.6	82	-0.03
33		4-Chloroaniline	0.420	0.450	-7.1	96	-0.01
34	C	Hexachlorobutadiene	0.329	0.324	1.5	93	-0.01
35		Caprolactam	0.126	0.121	4.0	90	-0.01
36	C	4-Chloro-3-methylphenol	0.412	0.421	-2.2	93	-0.01
37		2-Methylnaphthalene	0.740	0.765	-3.4	96	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	95	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.660	0.696	-5.5	95	-0.01
40	P	Hexachlorocyclopentadiene	0.192	0.237	-23.4	102	-0.01
41	S	2,4,6-Tribromophenol	0.322	0.313	2.8	86	-0.01
42	C	2,4,6-Trichlorophenol	0.416	0.419	-0.7	90	-0.01
43		2,4,5-Trichlorophenol	0.435	0.428	1.6	88	0.00
44	S	2-Fluorobiphenyl	1.235	1.305	-5.7	97	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.345	1.441	-7.1	97	-0.01
46		2-Chloronaphthalene	1.051	1.091	-3.8	95	-0.02
47		2-Nitroaniline	0.444	0.476	-7.2	94	-0.01
48		Acenaphthylene	1.629	1.719	-5.5	96	-0.02
49		Dimethylphthalate	1.458	1.561	-7.1	95	-0.01
50		2,6-Dinitrotoluene	0.319	0.327	-2.5	92	-0.01
51	C	Acenaphthene	1.065	1.110	-4.2	94	-0.01
52		3-Nitroaniline	0.306	0.316	-3.3	91	-0.01
53	P	2,4-Dinitrophenol	0.183	0.176	3.8	81	-0.01
54		Dibenzofuran	1.684	1.801	-6.9	96	-0.02
55	P	4-Nitrophenol	0.229	0.219	4.4	82	0.00
56		2,4-Dinitrotoluene	0.464	0.477	-2.8	91	-0.01
57		Fluorene	1.410	1.468	-4.1	94	-0.01
58		2,3,4,6-Tetrachlorophenol	0.466	0.453	2.8	88	-0.01
59		Diethylphthalate	1.529	1.574	-2.9	93	-0.01
60		4-Chlorophenyl-phenylether	0.799	0.811	-1.5	92	-0.01
61		4-Nitroaniline	0.308	0.308	0.0	81	-0.01
62		Azobenzene	1.475	1.610	-9.2	97	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	93	-0.01
64		4,6-Dinitro-2-methylphenol	0.139	0.140	-0.7	82	0.00
65	c	n-Nitrosodiphenylamine	0.541	0.558	-3.1	91	-0.02
66		4-Bromophenyl-phenylether	0.247	0.256	-3.6	91	-0.01
67		Hexachlorobenzene	0.283	0.283	0.0	89	-0.01
68		Atrazine	0.237	0.205	13.5	76	-0.01
69	C	Pentachlorophenol	0.147	0.131	10.9	75	0.00
70		Phenanthrene	1.031	1.063	-3.1	92	-0.01
71		Anthracene	1.047	1.076	-2.8	92	-0.01
72		Carbazole	0.937	0.950	-1.4	90	-0.01
73		Di-n-butylphthalate	1.152	1.175	-2.0	91	-0.01
74	C	Fluoranthene	1.361	1.416	-4.0	93	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	101	-0.02
76		Benzidine	0.499	0.465	6.8	89	-0.01
77		Pyrene	1.045	1.050	-0.5	93	-0.01
78	S	Terphenyl-d14	0.754	0.743	1.5	93	-0.01
79		Butylbenzylphthalate	0.420	0.417	0.7	96	-0.01
80		Benzo(a)anthracene	1.116	1.157	-3.7	98	-0.02
81		3,3'-Dichlorobenzidine	0.433	0.444	-2.5	96	-0.01
82		Chrysene	1.070	1.092	-2.1	98	-0.02
83		Bis(2-ethylhexyl)phthalate	0.584	0.619	-6.0	102	-0.02
84	c	Di-n-octyl phthalate	0.970	1.048	-8.0	102	-0.02
85		Indeno(1,2,3-cd)pyrene	1.360	1.483	-9.0	103	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	106	-0.03
87		Benzo(b)fluoranthene	1.091	1.092	-0.1	102	-0.02

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88	Benzo(k)fluoranthene	1.054	1.066	-1.1	101	-0.03
89 C	Benzo(a)pyrene	1.054	1.054	0.0	101	-0.03
90	Dibenzo(a,h)anthracene	1.117	1.143	-2.3	102	-0.05
91	Benzo(g,h,i)perylene	1.110	1.120	-0.9	103	-0.05

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