Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_G\Data\BG110316\
Data File : BG024683.D
Acq On : 4 Nov 2016 15:17
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Nov 04 15:52:38 2016
Quant Method : Z:\HPCHEM1\BNA_G\METHODS\8270-BG102516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 26 11:14:13 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	82	0.00
2		1,4-Dioxane	0.499	0.535	-7.2	92	0.00
3		Pyridine	1.493	1.628	-9.0	93	0.00
4		n-Nitrosodimethylamine	0.773	0.795	-2.8	86	0.00
5	S	2-Fluorophenol	1.220	1.262	-3.4	90	0.00
6		Aniline	2.121	2.288	-7.9	91	0.00
7	S	Phenol-d6	1.674	1.868	-11.6	94	0.00
8		2-Chlorophenol	1.241	1.393	-12.2	98	0.00
9		Benzaldehyde	1.034	1.045	-1.1	86	0.00
10	C	Phenol	1.725	1.963	-13.8	97	0.00
11		bis(2-Chloroethyl)ether	1.296	1.421	-9.6	96	0.00
12		1,3-Dichlorobenzene	1.536	1.583	-3.1	91	0.00
13	C	1,4-Dichlorobenzene	1.517	1.641	-8.2	93	0.00
14		1,2-Dichlorobenzene	1.459	1.533	-5.1	91	0.00
15		Benzyl Alcohol	1.557	1.700	-9.2	93	0.00
16		2,2'-oxybis(1-Chloropropane	2.405	2.448	-1.8	87	0.00
17		2-Methylphenol	1.143	1.281	-12.1	96	0.00
18		Hexachloroethane	0.583	0.636	-9.1	98	0.00
19	P	n-Nitroso-di-n-propylamine	1.234	1.341	-8.7	90	0.00
20		3+4-Methylphenols	1.535	1.699	-10.7	92	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	87	0.00
22		Acetophenone	0.514	0.512	0.4	91	0.00
23	S	Nitrobenzene-d5	0.439	0.450	-2.5	94	0.00
24		Nitrobenzene	0.489	0.497	-1.6	92	0.00
25		Isophorone	0.817	0.807	1.2	89	0.00
26	C	2-Nitrophenol	0.181	0.194	-7.2	98	0.00
27		2,4-Dimethylphenol	0.318	0.321	-0.9	90	0.00
28		bis(2-Chloroethoxy)methane	0.423	0.429	-1.4	95	0.00
29	C	2,4-Dichlorophenol	0.333	0.348	-4.5	94	0.00
30		1,2,4-Trichlorobenzene	0.395	0.404	-2.3	95	0.00
31		Naphthalene	0.998	1.016	-1.8	93	0.00
32		Benzoic acid	0.264	0.221	16.3	76	0.00
33		4-Chloroaniline	0.433	0.459	-6.0	92	0.00
34	C	Hexachlorobutadiene	0.309	0.316	-2.3	97	0.00
35		Caprolactam	0.122	0.122	0.0	89	0.00
36	C	4-Chloro-3-methylphenol	0.402	0.425	-5.7	93	0.00
37		2-Methylnaphthalene	0.728	0.765	-5.1	96	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	87	0.00
39		1,2,4,5-Tetrachlorobenzene	0.675	0.664	1.6	94	0.00
40	P	Hexachlorocyclopentadiene	0.380	0.328	13.7	80	0.00
41	S	2,4,6-Tribromophenol	0.299	0.317	-6.0	93	0.00
42	C	2,4,6-Trichlorophenol	0.405	0.432	-6.7	100	0.00
43		2,4,5-Trichlorophenol	0.438	0.456	-4.1	96	0.00
44	S	2-Fluorobiphenyl	1.203	1.216	-1.1	95	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.388	1.371	1.2	92	0.00
46		2-Chloronaphthalene	1.057	1.065	-0.8	95	0.00
47		2-Nitroaniline	0.433	0.461	-6.5	91	0.00
48		Acenaphthylene	1.621	1.656	-2.2	93	0.00
49		Dimethylphthalate	1.420	1.429	-0.6	90	0.00
50		2,6-Dinitrotoluene	0.303	0.315	-4.0	90	0.00
51	C	Acenaphthene	1.208	1.086	10.1	83	0.00
52		3-Nitroaniline	0.314	0.322	-2.5	90	0.00
53	P	2,4-Dinitrophenol	0.220	0.162	26.4#	63	0.00
54		Dibenzofuran	1.670	1.711	-2.5	94	0.00
55	P	4-Nitrophenol	0.273	0.251	8.1	81	0.00
56		2,4-Dinitrotoluene	0.440	0.483	-9.8	92	0.00
57		Fluorene	1.385	1.409	-1.7	91	0.00
58		2,3,4,6-Tetrachlorophenol	0.454	0.482	-6.2	93	0.00
59		Diethylphthalate	1.489	1.467	1.5	88	0.00
60		4-Chlorophenyl-phenylether	0.777	0.804	-3.5	95	0.00
61		4-Nitroaniline	0.343	0.365	-6.4	94	0.00
62		Azobenzene	1.485	1.481	0.3	89	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	83	0.00
64		4,6-Dinitro-2-methylphenol	0.138	0.131	5.1	82	0.00
65	c	n-Nitrosodiphenylamine	0.547	0.564	-3.1	92	0.00
66		4-Bromophenyl-phenylether	0.243	0.248	-2.1	92	0.00
67		Hexachlorobenzene	0.268	0.280	-4.5	92	0.00
68		Atrazine	0.235	0.232	1.3	86	0.00
69	C	Pentachlorophenol	0.177	0.166	6.2	79	0.00
70		Phenanthrene	1.019	1.046	-2.6	92	0.00
71		Anthracene	1.028	1.048	-1.9	90	0.00
72		Carbazole	0.938	0.945	-0.7	91	0.00
73		Di-n-butylphthalate	1.081	1.102	-1.9	90	0.00
74	C	Fluoranthene	1.289	1.332	-3.3	90	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	80	0.00
76		Benzidine	0.471	0.442	6.2	80	0.00
77		Pyrene	0.978	1.054	-7.8	89	0.00
78	S	Terphenyl-d14	0.669	0.699	-4.5	89	0.00
79		Butylbenzylphthalate	0.408	0.417	-2.2	87	0.00
80		Benzo(a)anthracene	1.099	1.133	-3.1	90	0.00
81		3,3'-Dichlorobenzidine	0.443	0.431	2.7	83	0.00
82		Chrysene	1.046	1.077	-3.0	88	0.00
83		Bis(2-ethylhexyl)phthalate	0.583	0.585	-0.3	86	0.00
84	c	Di-n-octyl phthalate	0.968	0.982	-1.4	86	-0.02
85		Indeno(1,2,3-cd)pyrene	1.444	1.393	3.5	86	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	79	-0.02
87		Benzo(b)fluoranthene	1.081	1.102	-1.9	87	0.00

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88	Benzo(k)fluoranthene	1.044	1.080	-3.4	87	-0.02
89 C	Benzo(a)pyrene	1.056	1.078	-2.1	86	-0.02
90	Dibenzo(a,h)anthracene	1.159	1.129	2.6	85	-0.02
91	Benzo(g,h,i)perylene	1.158	1.118	3.5	85	-0.02

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