Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF041816\
Data File : BF086275.D
Acq On : 18 Apr 2016 11:54
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Apr 19 01:02:10 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF041516.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 15 15:56:31 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	116	-0.01
2		1,4-Dioxane	0.673	0.694	-3.1	124	0.01
3		Pyridine	1.756	1.721	2.0	110	0.01
4		n-Nitrosodimethylamine	0.634	0.683	-7.7	128	0.01
5	S	2-Fluorophenol	1.221	1.098	10.1	115	0.00
6		Aniline	2.201	2.253	-2.4	121	0.00
7	S	Phenol-d6	1.527	1.485	2.8	116	0.00
8		2-Chlorophenol	1.381	1.305	5.5	114	0.00
9		Benzaldehyde	1.083	1.037	4.2	109	-0.01
10	C	Phenol	1.818	1.843	-1.4	113	0.00
11		bis(2-Chloroethyl)ether	1.368	1.250	8.6	112	0.00
12		1,3-Dichlorobenzene	1.465	1.407	4.0	120	0.00
13	C	1,4-Dichlorobenzene	1.396	1.474	-5.6	128	0.00
14		1,2-Dichlorobenzene	1.239	1.203	2.9	111	-0.01
15		Benzyl Alcohol	1.083	1.134	-4.7	117	0.00
16		2,2'-oxybis(1-Chloropropane	2.131	2.190	-2.8	119	0.00
17		2-Methylphenol	1.151	1.208	-5.0	122	0.00
18		Hexachloroethane	0.527	0.531	-0.8	116	-0.01
19	P	n-Nitroso-di-n-propylamine	0.977	0.863	11.7	119	0.00
20		3+4-Methylphenols	1.289	1.127	12.6	113	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	122	0.00
22		Acetophenone	0.421	0.374	11.2	113	0.00
23	S	Nitrobenzene-d5	0.347	0.338	2.6	126	0.00
24		Nitrobenzene	0.379	0.359	5.3	109	-0.01
25		Isophorone	0.688	0.669	2.8	125	0.00
26	C	2-Nitrophenol	0.165	0.176	-6.7	130	0.00
27		2,4-Dimethylphenol	0.334	0.319	4.5	117	-0.01
28		bis(2-Chloroethoxy)methane	0.405	0.365	9.9	122	0.00
29	C	2,4-Dichlorophenol	0.254	0.244	3.9	124	0.00
30		1,2,4-Trichlorobenzene	0.266	0.265	0.4	130	0.00
31		Naphthalene	0.893	0.893	0.0	128	0.00
32		Benzoic acid	0.213	0.233	-9.4	130	0.01
33		4-Chloroaniline	0.399	0.393	1.5	133	0.00
34	C	Hexachlorobutadiene	0.133	0.123	7.5	115	-0.01
35		Caprolactam	0.092	0.100	-8.7	129	0.00
36	C	4-Chloro-3-methylphenol	0.294	0.278	5.4	114	0.00
37		2-Methylnaphthalene	0.577	0.520	9.9	111	-0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	118	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.464	0.493	-6.2	133	0.00
40	P	Hexachlorocyclopentadiene	0.236	0.277	-17.4	144	0.00
41	S	2,4,6-Tribromophenol	0.161	0.158	1.9	122	-0.01
42	C	2,4,6-Trichlorophenol	0.394	0.405	-2.8	127	-0.01
43		2,4,5-Trichlorophenol	0.381	0.398	-4.5	122	0.00
44	S	2-Fluorobiphenyl	1.084	0.965	11.0	108	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.440	1.411	2.0	126	0.00
46		2-Chloronaphthalene	1.175	1.134	3.5	122	0.00
47		2-Nitroaniline	0.381	0.415	-8.9	136	0.00
48		Acenaphthylene	1.870	1.829	2.2	112	-0.01
49		Dimethylphthalate	1.454	1.434	1.4	119	-0.01
50		2,6-Dinitrotoluene	0.322	0.348	-8.1	122	0.00
51	C	Acenaphthene	1.117	1.124	-0.6	117	-0.01
52		3-Nitroaniline	0.391	0.362	7.4	125	0.00
53	P	2,4-Dinitrophenol	0.103	0.159	-54.4#	155#	0.00
54		Dibenzofuran	1.476	1.435	2.8	113	-0.01
55	P	4-Nitrophenol	0.306	0.357	-16.7	143	0.01
56		2,4-Dinitrotoluene	0.412	0.472	-14.6	130	0.00
57		Fluorene	1.227	0.972	20.8	112	-0.01
58		2,3,4,6-Tetrachlorophenol	0.276	0.266	3.6	121	-0.01
59		Diethylphthalate	1.501	1.437	4.3	122	0.00
60		4-Chlorophenyl-phenylether	0.506	0.490	3.2	133	0.00
61		4-Nitroaniline	0.338	0.405	-19.8	131	0.00
62		Azobenzene	1.469	1.546	-5.2	124	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	121	0.00
64		4,6-Dinitro-2-methylphenol	0.095	0.116	-22.1	139	0.00
65	c	n-Nitrosodiphenylamine	0.612	0.567	7.4	122	0.00
66		4-Bromophenyl-phenylether	0.186	0.188	-1.1	126	0.00
67		Hexachlorobenzene	0.195	0.182	6.7	118	0.00
68		Atrazine	0.188	0.170	9.6	115	0.00
69	C	Pentachlorophenol	0.118	0.135	-14.4	130	0.00
70		Phenanthrene	0.955	0.877	8.2	112	0.00
71		Anthracene	1.043	1.011	3.1	131	0.00
72		Carbazole	0.908	0.856	5.7	122	0.00
73		Di-n-butylphthalate	1.228	1.211	1.4	128	0.00
74	C	Fluoranthene	0.978	0.914	6.5	129	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	118	0.00
76		Benzidine	0.776	0.851	-9.7	117	0.00
77		Pyrene	1.534	1.402	8.6	126	0.00
78	S	Terphenyl-d14	0.795	0.758	4.7	131	0.00
79		Butylbenzylphthalate	0.709	0.738	-4.1	115	-0.01
80		Benzo(a)anthracene	1.098	1.001	8.8	122	0.00
81		3,3'-Dichlorobenzidine	0.636	0.683	-7.4	133	0.00
82		Chrysene	1.171	1.079	7.9	120	0.00
83		Bis(2-ethylhexyl)phthalate	0.765	0.780	-2.0	123	-0.01
84	c	Di-n-octyl phthalate	1.475	1.624	-10.1	122	0.00
85		Indeno(1,2,3-cd)pyrene	0.962	1.039	-8.0	135	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	133	0.00
87		Benzo(b)fluoranthene	1.234	1.150	6.8	136	0.00

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88	Benzo(k)fluoranthene	1.125	1.044	7.2	117	-0.01
89 C	Benzo(a)pyrene	1.069	1.059	0.9	131	0.00
90	Dibenzo(a,h)anthracene	0.945	0.923	2.3	134	0.00
91	Benzo(g,h,i)perylene	1.004	0.982	2.2	135	0.00

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