Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071415\
Data File : BF080492.D
Acq On : 15 Jul 2015 14:57
Operator : TP/UM
Sample : SSTDCCC040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 15 16:17:47 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jul 15 03:58:42 2015
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	143	0.00
2		1,4-Dioxane	0.493	0.512	-3.9	159#	0.00
3		Pyridine	1.064	1.363	-28.1#	162#	-0.03
4		n-Nitrosodimethylamine	0.679	0.807	-18.9	165#	-0.01
5	S	2-Fluorophenol	1.224	1.226	-0.2	145	0.00
6		Aniline	2.174	2.169	0.2	139	0.00
7	S	Phenol-d6	1.497	1.550	-3.5	149	-0.01
8		2-Chlorophenol	1.284	1.270	1.1	142	0.00
9		Benzaldehyde	0.914	0.953	-4.3	145	0.00
10	C	Phenol	1.726	1.664	3.6	138	0.00
11		bis(2-Chloroethyl)ether	1.312	1.271	3.1	136	0.00
12		1,3-Dichlorobenzene	1.497	1.529	-2.1	147	0.00
13	C	1,4-Dichlorobenzene	1.627	1.661	-2.1	151#	0.00
14		1,2-Dichlorobenzene	1.573	1.543	1.9	144	0.00
15		Benzyl Alcohol	1.072	1.065	0.7	139	0.00
16		2,2'-oxybis(1-Chloropropane	2.127	2.183	-2.6	149	0.00
17		2-Methylphenol	1.039	1.054	-1.4	146	0.00
18		Hexachloroethane	0.536	0.546	-1.9	147	0.00
19	P	n-Nitroso-di-n-propylamine	1.014	1.008	0.6	135	0.00
20		3+4-Methylphenols	1.472	1.472	0.0	137	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	145	0.00
22		Acetophenone	0.469	0.472	-0.6	151#	-0.01
23	S	Nitrobenzene-d5	0.365	0.365	0.0	146	0.00
24		Nitrobenzene	0.403	0.387	4.0	140	0.00
25		Isophorone	0.675	0.701	-3.9	154#	0.00
26	C	2-Nitrophenol	0.175	0.172	1.7	140	0.00
27		2,4-Dimethylphenol	0.320	0.321	-0.3	142	0.00
28		bis(2-Chloroethoxy)methane	0.379	0.376	0.8	152#	0.00
29	C	2,4-Dichlorophenol	0.279	0.280	-0.4	154#	0.00
30		1,2,4-Trichlorobenzene	0.348	0.333	4.3	143	0.00
31		Naphthalene	1.035	0.996	3.8	146	0.00
32		Benzoic acid	0.152	0.172	-13.2	167#	0.00
33		4-Chloroaniline	0.412	0.413	-0.2	143	0.00
34	C	Hexachlorobutadiene	0.176	0.169	4.0	138	0.00
35		Caprolactam	0.072	0.076	-5.6	156#	-0.01
36	C	4-Chloro-3-methylphenol	0.273	0.276	-1.1	153#	-0.01
37		2-Methylnaphthalene	0.705	0.644	8.7	133	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	139	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.664	0.641	3.5	147	0.00
40	P	Hexachlorocyclopentadiene	0.299	0.184	38.5#	92	0.00
41	S	2,4,6-Tribromophenol	0.189	0.190	-0.5	134	0.00
42	C	2,4,6-Trichlorophenol	0.413	0.405	1.9	133	0.00
43		2,4,5-Trichlorophenol	0.416	0.417	-0.2	139	-0.01
44	S	2-Fluorobiphenyl	1.460	1.406	3.7	143	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.722	1.679	2.5	148	0.00
46		2-Chloronaphthalene	1.414	1.350	4.5	134	0.00
47		2-Nitroaniline	0.408	0.423	-3.7	149	0.00
48		Acenaphthylene	2.138	2.113	1.2	143	0.00
49		Dimethylphthalate	1.583	1.599	-1.0	150#	0.00
50		2,6-Dinitrotoluene	0.341	0.350	-2.6	148	0.00
51	C	Acenaphthene	1.301	1.242	4.5	137	0.00
52		3-Nitroaniline	0.367	0.381	-3.8	150#	0.00
53	P	2,4-Dinitrophenol	0.117	0.068	41.9#	76	0.00
54		Dibenzofuran	1.834	1.749	4.6	142	0.00
55	P	4-Nitrophenol	0.241	0.248	-2.9	137	0.00
56		2,4-Dinitrotoluene	0.386	0.394	-2.1	149	-0.01
57		Fluorene	1.475	1.460	1.0	141	0.00
58		2,3,4,6-Tetrachlorophenol	0.337	0.364	-8.0	157#	0.00
59		Diethylphthalate	1.567	1.573	-0.4	149	0.00
60		4-Chlorophenyl-phenylether	0.692	0.696	-0.6	150	0.00
61		4-Nitroaniline	0.333	0.365	-9.6	152#	0.00
62		Azobenzene	1.497	1.556	-3.9	149	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	141	0.00
64		4,6-Dinitro-2-methylphenol	0.101	0.060	40.6#	86	0.00
65	c	n-Nitrosodiphenylamine	0.650	0.609	6.3	137	0.00
66		4-Bromophenyl-phenylether	0.199	0.195	2.0	151#	0.00
67		Hexachlorobenzene	0.217	0.197	9.2	131	0.00
68		Atrazine	0.189	0.200	-5.8	159#	0.00
69	C	Pentachlorophenol	0.100	0.103	-3.0	149	0.00
70		Phenanthrene	1.162	1.057	9.0	137	0.00
71		Anthracene	1.093	1.130	-3.4	151#	0.00
72		Carbazole	0.990	1.015	-2.5	154#	0.00
73		Di-n-butylphthalate	1.051	1.214	-15.5	178#	0.02
74	C	Fluoranthene	1.146	1.202	-4.9	158#	0.05
75	I	Chrysene-d12	1.000	1.000	0.0	157#	0.14
76		Benzidine	0.593	0.551	7.1	138	0.06
77		Pyrene	1.177	1.166	0.9	158#	0.06
78	S	Terphenyl-d14	0.830	0.834	-0.5	157#	0.07
79		Butylbenzylphthalate	0.430	0.474	-10.2	171#	0.10
80		Benzo(a)anthracene	1.166	1.152	1.2	167#	0.14
81		3,3'-Dichlorobenzidine	0.368	0.377	-2.4	163#	0.14
82		Chrysene	1.114	1.063	4.6	151#	0.14
83		Bis(2-ethylhexyl)phthalate	0.625	0.711	-13.8	179#	0.14
84	c	Di-n-octyl phthalate	1.122	1.212	-8.0	170#	0.17
85		Indeno(1,2,3-cd)pyrene	1.780	1.223	31.3#	111	0.22
86	I	Perylene-d12	1.000	1.000	0.0	140	0.21
87		Benzo(b)fluoranthene	1.143	1.119	2.1	152#	0.19

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88	Benzo(k)fluoranthene	1.104	1.116	-1.1	138	0.19
89 C	Benzo(a)pyrene	1.125	1.051	6.6	135	0.19
90	Dibenzo(a,h)anthracene	1.276	1.028	19.4	113	0.22
91	Benzo(g,h,i)perylene	1.336	0.992	25.7#	104	0.21

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