Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF020615\
Data File : BF077220.D
Acq On : 7 Feb 2015 1:31
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : SSTDCCC040EC

Quant Time: Feb 07 02:37:24 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Feb 07 02:27:13 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	130	-0.04
2		1,4-Dioxane	0.521	0.441	15.4	116	-0.03
3		Pyridine	1.348	1.401	-3.9	112	-0.04
4		n-Nitrosodimethylamine	0.668	0.644	3.6	124	-0.03
5	S	2-Fluorophenol	1.216	1.187	2.4	121	-0.05
6		Aniline	2.124	2.117	0.3	122	-0.04
7	S	Phenol-d6	1.535	1.500	2.3	121	-0.03
8		2-Chlorophenol	1.402	1.384	1.3	121	-0.04
9		Benzaldehyde	0.894	0.837	6.4	139	-0.04
10	C	Phenol	1.629	1.574	3.4	115	-0.04
11		bis(2-Chloroethyl)ether	1.275	1.294	-1.5	128	-0.04
12		1,3-Dichlorobenzene	1.595	1.619	-1.5	128	-0.04
13	C	1,4-Dichlorobenzene	1.692	1.665	1.6	125	-0.05
14		1,2-Dichlorobenzene	1.559	1.543	1.0	122	-0.04
15		Benzyl Alcohol	1.242	1.299	-4.6	126	-0.04
16		2,2'-oxybis(1-Chloropropane	1.714	1.723	-0.5	126	-0.04
17		2-Methylphenol	1.146	1.146	0.0	121	-0.04
18		Hexachloroethane	0.588	0.616	-4.8	129	-0.04
19	P	n-Nitroso-di-n-propylamine	1.074	1.082	-0.7	123	-0.03
20		3+4-Methylphenols	1.517	1.394	8.1	114	-0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	131	-0.04
22		Acetophenone	0.490	0.488	0.4	123	-0.04
23	S	Nitrobenzene-d5	0.361	0.374	-3.6	127	-0.04
24		Nitrobenzene	0.368	0.363	1.4	119	-0.04
25		Isophorone	0.657	0.673	-2.4	125	-0.03
26	C	2-Nitrophenol	0.172	0.172	0.0	115	-0.04
27		2,4-Dimethylphenol	0.284	0.290	-2.1	122	-0.03
28		bis(2-Chloroethoxy)methane	0.374	0.386	-3.2	123	-0.03
29	C	2,4-Dichlorophenol	0.265	0.249	6.0	113	-0.03
30		1,2,4-Trichlorobenzene	0.294	0.301	-2.4	121	-0.04
31		Naphthalene	1.030	1.033	-0.3	122	-0.04
32		Benzoic acid	0.158	0.085	46.2#	65	0.00
33		4-Chloroaniline	0.416	0.394	5.3	116	-0.03
34	C	Hexachlorobutadiene	0.175	0.182	-4.0	126	-0.04
35		Caprolactam	0.078	0.074	5.1	110	-0.02
36	C	4-Chloro-3-methylphenol	0.303	0.295	2.6	118	-0.04
37		2-Methylnaphthalene	0.703	0.709	-0.9	125	-0.04
38	I	Acenaphthene-d10	1.000	1.000	0.0	124	-0.05
39		1,2,4,5-Tetrachlorobenzene	0.494	0.500	-1.2	122	-0.04
40	P	Hexachlorocyclopentadiene	0.202	0.241	-19.3	120	-0.04
41	S	2,4,6-Tribromophenol	0.162	0.168	-3.7	119	-0.03
42	C	2,4,6-Trichlorophenol	0.360	0.361	-0.3	115	-0.04
43		2,4,5-Trichlorophenol	0.367	0.323	12.0	101	-0.04
44	S	2-Fluorobiphenyl	1.314	1.357	-3.3	123	-0.04

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.483	1.541	-3.9	121	-0.04
46		2-Chloronaphthalene	1.220	1.256	-3.0	124	-0.04
47		2-Nitroaniline	0.370	0.403	-8.9	124	-0.03
48		Acenaphthylene	2.058	2.116	-2.8	121	-0.04
49		Dimethylphthalate	1.418	1.470	-3.7	125	-0.03
50		2,6-Dinitrotoluene	0.283	0.304	-7.4	120	-0.04
51	C	Acenaphthene	1.168	1.191	-2.0	121	-0.04
52		3-Nitroaniline	0.339	0.343	-1.2	113	-0.03
53	P	2,4-Dinitrophenol	0.095	0.085	10.5	99	-0.02
54		Dibenzofuran	1.701	1.719	-1.1	121	-0.04
55	P	4-Nitrophenol	0.222	0.168	24.3	87	-0.02
56		2,4-Dinitrotoluene	0.384	0.409	-6.5	119	-0.04
57		Fluorene	1.441	1.461	-1.4	121	-0.04
58		2,3,4,6-Tetrachlorophenol	0.268	0.261	2.6	112	-0.04
59		Diethylphthalate	1.433	1.532	-6.9	127	-0.04
60		4-Chlorophenyl-phenylether	0.590	0.606	-2.7	123	-0.03
61		4-Nitroaniline	0.328	0.341	-4.0	117	-0.03
62		Azobenzene	1.494	1.553	-3.9	124	-0.03
63	I	Phenanthrene-d10	1.000	1.000	0.0	131	-0.03
64		4,6-Dinitro-2-methylphenol	0.104	0.120	-15.4	119	-0.03
65	c	n-Nitrosodiphenylamine	0.713	0.696	2.4	118	-0.03
66		4-Bromophenyl-phenylether	0.203	0.204	-0.5	120	-0.03
67		Hexachlorobenzene	0.214	0.216	-0.9	126	-0.03
68		Atrazine	0.216	0.227	-5.1	120	-0.03
69	C	Pentachlorophenol	0.082	0.093	-13.4	134	-0.03
70		Phenanthrene	1.247	1.200	3.8	120	-0.03
71		Anthracene	1.216	1.192	2.0	123	-0.03
72		Carbazole	1.180	1.177	0.3	122	-0.03
73		Di-n-butylphthalate	1.365	1.462	-7.1	129	-0.03
74	C	Fluoranthene	1.278	1.294	-1.3	122	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	135	-0.03
76		Benzidine	0.640	0.607	5.2	129	-0.02
77		Pyrene	1.371	1.403	-2.3	126	-0.02
78	S	Terphenyl-d14	0.869	0.866	0.3	125	-0.03
79		Butylbenzylphthalate	0.621	0.657	-5.8	126	-0.03
80		Benzo(a)anthracene	1.255	1.256	-0.1	125	-0.03
81		3,3'-Dichlorobenzidine	0.399	0.423	-6.0	133	-0.02
82		Chrysene	1.144	1.148	-0.3	125	-0.03
83		Bis(2-ethylhexyl)phthalate	0.859	0.937	-9.1	137	-0.03
84	c	Di-n-octyl phthalate	1.491	1.670	-12.0	138	-0.03
85		Indeno(1,2,3-cd)pyrene	1.213	1.338	-10.3	136	-0.10
86	I	Perylene-d12	1.000	1.000	0.0	134	-0.05
87		Benzo(b)fluoranthene	1.347	1.277	5.2	133	-0.04

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88	Benzo(k)fluoranthene	1.177	1.336	-13.5	133	-0.03
89 C	Benzo(a)pyrene	1.188	1.189	-0.1	127	-0.04
90	Dibenzo(a,h)anthracene	1.090	1.151	-5.6	133	-0.10
91	Benzo(g,h,i)perylene	1.084	1.144	-5.5	133	-0.10

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