Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF012215\
Data File : BF077010.D
Acq On : 22 Jan 2015 12:38
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 22 23:55:04 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jan 21 14:45:53 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	93	0.00
2		1,4-Dioxane	0.521	0.419	19.6	79	0.00
3		Pyridine	1.348	1.334	1.0	77	-0.01
4		n-Nitrosodimethylamine	0.668	0.684	-2.4	95	-0.01
5	S	2-Fluorophenol	1.216	1.222	-0.5	89	-0.01
6		Aniline	2.124	2.149	-1.2	88	0.00
7	S	Phenol-d6	1.535	1.518	1.1	88	-0.01
8		2-Chlorophenol	1.402	1.397	0.4	87	0.00
9		Benzaldehyde	0.894	0.919	-2.8	109	0.00
10	C	Phenol	1.629	1.664	-2.1	87	-0.01
11		bis(2-Chloroethyl)ether	1.275	1.303	-2.2	92	0.00
12		1,3-Dichlorobenzene	1.595	1.588	0.4	90	-0.01
13	C	1,4-Dichlorobenzene	1.692	1.652	2.4	89	-0.01
14		1,2-Dichlorobenzene	1.559	1.568	-0.6	89	0.00
15		Benzyl Alcohol	1.242	1.262	-1.6	88	-0.01
16		2,2'-oxybis(1-Chloropropane	1.714	1.680	2.0	88	0.00
17		2-Methylphenol	1.146	1.146	0.0	87	-0.01
18		Hexachloroethane	0.588	0.579	1.5	87	-0.01
19	P	n-Nitroso-di-n-propylamine	1.074	1.057	1.6	86	0.00
20		3+4-Methylphenols	1.517	1.503	0.9	88	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	91	0.00
22		Acetophenone	0.490	0.507	-3.5	89	0.00
23	S	Nitrobenzene-d5	0.361	0.376	-4.2	89	-0.01
24		Nitrobenzene	0.368	0.373	-1.4	85	0.00
25		Isophorone	0.657	0.680	-3.5	88	-0.01
26	C	2-Nitrophenol	0.172	0.183	-6.4	85	-0.01
27		2,4-Dimethylphenol	0.284	0.297	-4.6	87	-0.01
28		bis(2-Chloroethoxy)methane	0.374	0.392	-4.8	87	0.00
29	C	2,4-Dichlorophenol	0.265	0.275	-3.8	87	-0.01
30		1,2,4-Trichlorobenzene	0.294	0.303	-3.1	85	0.00
31		Naphthalene	1.030	1.063	-3.2	88	0.00
32		Benzoic acid	0.158	0.148	6.3	79	-0.02
33		4-Chloroaniline	0.416	0.435	-4.6	90	-0.01
34	C	Hexachlorobutadiene	0.175	0.177	-1.1	85	0.00
35		Caprolactam	0.078	0.080	-2.6	84	-0.01
36	C	4-Chloro-3-methylphenol	0.303	0.308	-1.7	86	-0.01
37		2-Methylnaphthalene	0.703	0.716	-1.8	88	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	88	0.00
39		1,2,4,5-Tetrachlorobenzene	0.494	0.502	-1.6	87	0.00
40	P	Hexachlorocyclopentadiene	0.202	0.264	-30.7#	93	0.00
41	S	2,4,6-Tribromophenol	0.162	0.167	-3.1	83	-0.01
42	C	2,4,6-Trichlorophenol	0.360	0.371	-3.1	83	-0.01
43		2,4,5-Trichlorophenol	0.367	0.387	-5.4	86	0.00
44	S	2-Fluorobiphenyl	1.314	1.347	-2.5	86	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.483	1.548	-4.4	86	-0.01
46		2-Chloronaphthalene	1.220	1.269	-4.0	88	-0.01
47		2-Nitroaniline	0.370	0.399	-7.8	87	-0.01
48		Acenaphthylene	2.058	2.102	-2.1	85	-0.01
49		Dimethylphthalate	1.418	1.458	-2.8	87	0.00
50		2,6-Dinitrotoluene	0.283	0.306	-8.1	86	-0.01
51	C	Acenaphthene	1.168	1.188	-1.7	85	-0.01
52		3-Nitroaniline	0.339	0.353	-4.1	82	0.00
53	P	2,4-Dinitrophenol	0.095	0.107	-12.6	88	-0.01
54		Dibenzofuran	1.701	1.740	-2.3	87	0.00
55	P	4-Nitrophenol	0.222	0.228	-2.7	83	-0.01
56		2,4-Dinitrotoluene	0.384	0.415	-8.1	85	0.00
57		Fluorene	1.441	1.460	-1.3	86	0.00
58		2,3,4,6-Tetrachlorophenol	0.268	0.293	-9.3	89	0.00
59		Diethylphthalate	1.433	1.478	-3.1	86	0.00
60		4-Chlorophenyl-phenylether	0.590	0.590	0.0	85	0.00
61		4-Nitroaniline	0.328	0.356	-8.5	86	0.00
62		Azobenzene	1.494	1.552	-3.9	87	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	90	0.00
64		4,6-Dinitro-2-methylphenol	0.104	0.124	-19.2	84	0.00
65	c	n-Nitrosodiphenylamine	0.713	0.735	-3.1	85	-0.01
66		4-Bromophenyl-phenylether	0.203	0.213	-4.9	86	0.00
67		Hexachlorobenzene	0.214	0.222	-3.7	89	-0.01
68		Atrazine	0.216	0.228	-5.6	83	0.00
69	C	Pentachlorophenol	0.082	0.096	-17.1	95	-0.01
70		Phenanthrene	1.247	1.258	-0.9	86	0.00
71		Anthracene	1.216	1.232	-1.3	87	0.00
72		Carbazole	1.180	1.250	-5.9	89	0.00
73		Di-n-butylphthalate	1.365	1.470	-7.7	89	-0.01
74	C	Fluoranthene	1.278	1.313	-2.7	85	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	93	-0.01
76		Benzidine	0.640	0.712	-11.2	104	0.00
77		Pyrene	1.371	1.367	0.3	85	0.00
78	S	Terphenyl-d14	0.869	0.860	1.0	86	-0.01
79		Butylbenzylphthalate	0.621	0.650	-4.7	86	0.00
80		Benzo(a)anthracene	1.255	1.275	-1.6	87	-0.01
81		3,3'-Dichlorobenzidine	0.399	0.407	-2.0	88	0.00
82		Chrysene	1.144	1.131	1.1	85	0.00
83		Bis(2-ethylhexyl)phthalate	0.859	0.917	-6.8	92	0.00
84	c	Di-n-octyl phthalate	1.491	1.578	-5.8	90	-0.01
85		Indeno(1,2,3-cd)pyrene	1.213	1.244	-2.6	87	-0.03
86	I	Perylene-d12	1.000	1.000	0.0	91	-0.02
87		Benzo(b)fluoranthene	1.347	1.400	-3.9	98	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.177	1.161	1.4	78	0.02
89 C	Benzo(a)pyrene	1.188	1.193	-0.4	86	-0.02
90	Dibenzo(a,h)anthracene	1.090	1.131	-3.8	88	-0.03
91	Benzo(g,h,i)perylene	1.084	1.140	-5.2	90	-0.03

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