Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF011915\
Data File : BF076942.D
Acq On : 20 Jan 2015 00:00
Operator : IZ / TP
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jan 20 10:07:24 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jan 19 23:54:01 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	109	-0.05
2		1,4-Dioxane	0.521	0.540	-3.6	119	-0.05
3		Pyridine	1.348	1.605	-19.1	108	-0.05
4		n-Nitrosodimethylamine	0.668	0.745	-11.5	121	-0.05
5	S	2-Fluorophenol	1.216	1.290	-6.1	110	-0.06
6		Aniline	2.124	2.269	-6.8	109	-0.05
7	S	Phenol-d6	1.535	1.630	-6.2	110	-0.05
8		2-Chlorophenol	1.402	1.489	-6.2	109	-0.05
9		Benzaldehyde	0.894	0.826	7.6	115	-0.05
10	C	Phenol	1.629	1.772	-8.8	108	-0.05
11		bis(2-Chloroethyl)ether	1.275	1.421	-11.5	118	-0.05
12		1,3-Dichlorobenzene	1.595	1.757	-10.2	117	-0.05
13	C	1,4-Dichlorobenzene	1.692	1.789	-5.7	112	-0.05
14		1,2-Dichlorobenzene	1.559	1.702	-9.2	113	-0.05
15		Benzyl Alcohol	1.242	1.349	-8.6	110	-0.03
16		2,2'-oxybis(1-Chloropropane	1.714	1.807	-5.4	111	-0.05
17		2-Methylphenol	1.146	1.244	-8.6	110	-0.05
18		Hexachloroethane	0.588	0.645	-9.7	113	-0.05
19	P	n-Nitroso-di-n-propylamine	1.074	1.142	-6.3	109	-0.05
20		3+4-Methylphenols	1.517	1.639	-8.0	112	-0.05
21	I	Naphthalene-d8	1.000	1.000	0.0	113	-0.05
22		Acetophenone	0.490	0.523	-6.7	113	-0.05
23	S	Nitrobenzene-d5	0.361	0.381	-5.5	111	-0.03
24		Nitrobenzene	0.368	0.386	-4.9	109	-0.05
25		Isophorone	0.657	0.684	-4.1	109	-0.05
26	C	2-Nitrophenol	0.172	0.189	-9.9	109	-0.04
27		2,4-Dimethylphenol	0.284	0.292	-2.8	106	-0.03
28		bis(2-Chloroethoxy)methane	0.374	0.393	-5.1	108	-0.05
29	C	2,4-Dichlorophenol	0.265	0.289	-9.1	112	-0.05
30		1,2,4-Trichlorobenzene	0.294	0.318	-8.2	110	-0.05
31		Naphthalene	1.030	1.065	-3.4	109	-0.05
32		Benzoic acid	0.158	0.180	-13.9	119	-0.03
33		4-Chloroaniline	0.416	0.436	-4.8	111	-0.03
34	C	Hexachlorobutadiene	0.175	0.180	-2.9	108	-0.05
35		Caprolactam	0.078	0.083	-6.4	106	-0.05
36	C	4-Chloro-3-methylphenol	0.303	0.309	-2.0	107	-0.05
37		2-Methylnaphthalene	0.703	0.736	-4.7	112	-0.05
38	I	Acenaphthene-d10	1.000	1.000	0.0	107	-0.06
39		1,2,4,5-Tetrachlorobenzene	0.494	0.520	-5.3	110	-0.06
40	P	Hexachlorocyclopentadiene	0.202	0.261	-29.2#	112	-0.06
41	S	2,4,6-Tribromophenol	0.162	0.177	-9.3	108	-0.03
42	C	2,4,6-Trichlorophenol	0.360	0.379	-5.3	104	-0.05
43		2,4,5-Trichlorophenol	0.367	0.396	-7.9	107	-0.06
44	S	2-Fluorobiphenyl	1.314	1.417	-7.8	111	-0.06

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.483	1.564	-5.5	106	-0.06
46		2-Chloronaphthalene	1.220	1.303	-6.8	111	-0.06
47		2-Nitroaniline	0.370	0.408	-10.3	108	-0.05
48		Acenaphthylene	2.058	2.186	-6.2	108	-0.06
49		Dimethylphthalate	1.418	1.466	-3.4	107	-0.06
50		2,6-Dinitrotoluene	0.283	0.325	-14.8	111	-0.05
51	C	Acenaphthene	1.168	1.245	-6.6	109	-0.06
52		3-Nitroaniline	0.339	0.370	-9.1	105	-0.06
53	P	2,4-Dinitrophenol	0.095	0.110	-15.8	110	-0.06
54		Dibenzofuran	1.701	1.809	-6.3	110	-0.05
55	P	4-Nitrophenol	0.222	0.222	0.0	99	-0.05
56		2,4-Dinitrotoluene	0.384	0.433	-12.8	109	-0.05
57		Fluorene	1.441	1.513	-5.0	109	-0.05
58		2,3,4,6-Tetrachlorophenol	0.268	0.287	-7.1	106	-0.06
59		Diethylphthalate	1.433	1.568	-9.4	112	-0.05
60		4-Chlorophenyl-phenylether	0.590	0.631	-6.9	111	-0.03
61		4-Nitroaniline	0.328	0.366	-11.6	108	-0.04
62		Azobenzene	1.494	1.583	-6.0	109	-0.05
63	I	Phenanthrene-d10	1.000	1.000	0.0	110	-0.05
64		4,6-Dinitro-2-methylphenol	0.104	0.130	-25.0#	108	-0.04
65	c	n-Nitrosodiphenylamine	0.713	0.756	-6.0	107	-0.03
66		4-Bromophenyl-phenylether	0.203	0.221	-8.9	109	-0.03
67		Hexachlorobenzene	0.214	0.230	-7.5	113	-0.03
68		Atrazine	0.216	0.240	-11.1	107	-0.03
69	C	Pentachlorophenol	0.082	0.096	-17.1	116	-0.03
70		Phenanthrene	1.247	1.290	-3.4	108	-0.03
71		Anthracene	1.216	1.263	-3.9	109	-0.05
72		Carbazole	1.180	1.239	-5.0	108	-0.03
73		Di-n-butylphthalate	1.365	1.455	-6.6	108	-0.03
74	C	Fluoranthene	1.278	1.377	-7.7	109	-0.03
75	I	Chrysene-d12	1.000	1.000	0.0	116	-0.02
76		Benzidine	0.640	0.619	3.3	112	-0.02
77		Pyrene	1.371	1.458	-6.3	113	-0.03
78	S	Terphenyl-d14	0.869	0.887	-2.1	110	-0.03
79		Butylbenzylphthalate	0.621	0.663	-6.8	109	-0.03
80		Benzo(a)anthracene	1.255	1.268	-1.0	108	-0.02
81		3,3'-Dichlorobenzidine	0.399	0.415	-4.0	112	-0.03
82		Chrysene	1.144	1.154	-0.9	107	-0.03
83		Bis(2-ethylhexyl)phthalate	0.859	0.937	-9.1	117	-0.03
84	c	Di-n-octyl phthalate	1.491	1.568	-5.2	111	-0.05
85		Indeno(1,2,3-cd)pyrene	1.213	1.232	-1.6	107	-0.13
86	I	Perylene-d12	1.000	1.000	0.0	109	-0.08
87		Benzo(b)fluoranthene	1.347	1.442	-7.1	122	-0.06

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.177	1.268	-7.7	103	-0.07
89 C	Benzo(a)pyrene	1.188	1.280	-7.7	111	-0.08
90	Dibenzo(a,h)anthracene	1.090	1.222	-12.1	115	-0.14
91	Benzo(g,h,i)perylene	1.084	1.241	-14.5	117	-0.14

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