Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF020915\
Data File : BF077222.D
Acq On : 9 Feb 2015 13:04
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Feb 09 15:51:38 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF011415.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 09 15:41:22 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	77	0.00
2		1,4-Dioxane	0.521	0.507	2.7	79	0.00
3		Pyridine	1.348	1.604	-19.0	76	0.00
4		n-Nitrosodimethylamine	0.668	0.733	-9.7	84	0.00
5	S	2-Fluorophenol	1.216	1.255	-3.2	76	0.00
6		Aniline	2.124	2.098	1.2	71	0.00
7	S	Phenol-d6	1.535	1.523	0.8	73	0.00
8		2-Chlorophenol	1.402	1.450	-3.4	75	0.00
9		Benzaldehyde	0.894	0.786	12.1	77	0.00
10	C	Phenol	1.629	1.507	7.5	65	0.00
11		bis(2-Chloroethyl)ether	1.275	1.334	-4.6	78	0.00
12		1,3-Dichlorobenzene	1.595	1.694	-6.2	79	0.00
13	C	1,4-Dichlorobenzene	1.692	1.715	-1.4	76	0.00
14		1,2-Dichlorobenzene	1.559	1.612	-3.4	76	0.00
15		Benzyl Alcohol	1.242	1.275	-2.7	73	0.00
16		2,2'-oxybis(1-Chloropropane	1.714	1.878	-9.6	81	0.00
17		2-Methylphenol	1.146	1.164	-1.6	73	0.00
18		Hexachloroethane	0.588	0.635	-8.0	79	0.00
19	P	n-Nitroso-di-n-propylamine	1.074	1.113	-3.6	75	0.00
20		3+4-Methylphenols	1.517	1.310	13.6	63	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	74	0.00
22		Acetophenone	0.490	0.526	-7.3	75	0.00
23	S	Nitrobenzene-d5	0.361	0.389	-7.8	74	0.00
24		Nitrobenzene	0.368	0.386	-4.9	71	0.00
25		Isophorone	0.657	0.689	-4.9	72	0.00
26	C	2-Nitrophenol	0.172	0.175	-1.7	66	0.00
27		2,4-Dimethylphenol	0.284	0.302	-6.3	72	0.00
28		bis(2-Chloroethoxy)methane	0.374	0.410	-9.6	74	0.00
29	C	2,4-Dichlorophenol	0.265	0.257	3.0	65	0.00
30		1,2,4-Trichlorobenzene	0.294	0.321	-9.2	73	0.00
31		Naphthalene	1.030	1.113	-8.1	74	0.00
32		Benzoic acid	0.158	0.087	44.9#	38#	0.00
33		4-Chloroaniline	0.416	0.412	1.0	69	0.00
34	C	Hexachlorobutadiene	0.175	0.190	-8.6	74	0.00
35		Caprolactam	0.078	0.073	6.4	62	0.00
36	C	4-Chloro-3-methylphenol	0.303	0.298	1.7	67	0.00
37		2-Methylnaphthalene	0.703	0.760	-8.1	76	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	68	0.00
39		1,2,4,5-Tetrachlorobenzene	0.494	0.541	-9.5	73	0.00
40	P	Hexachlorocyclopentadiene	0.202	0.236	-16.8	64	0.00
41	S	2,4,6-Tribromophenol	0.162	0.167	-3.1	65	0.00
42	C	2,4,6-Trichlorophenol	0.360	0.366	-1.7	64	0.00
43		2,4,5-Trichlorophenol	0.367	0.301	18.0	52	0.00
44	S	2-Fluorobiphenyl	1.314	1.483	-12.9	73	0.00

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
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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)
45		1,1'-Biphenyl	1.483	1.651	-11.3	71	0.00
46		2-Chloronaphthalene	1.220	1.342	-10.0	72	0.00
47		2-Nitroaniline	0.370	0.392	-5.9	66	0.00
48		Acenaphthylene	2.058	2.263	-10.0	71	0.00
49		Dimethylphthalate	1.418	1.568	-10.6	73	0.00
50		2,6-Dinitrotoluene	0.283	0.323	-14.1	70	0.00
51	C	Acenaphthene	1.168	1.266	-8.4	71	0.00
52		3-Nitroaniline	0.339	0.351	-3.5	63	0.00
53	P	2,4-Dinitrophenol	0.095	0.102	-7.4	65	0.00
54		Dibenzofuran	1.701	1.852	-8.9	71	0.00
55	P	4-Nitrophenol	0.222	0.187	15.8	53	0.00
56		2,4-Dinitrotoluene	0.384	0.427	-11.2	68	0.00
57		Fluorene	1.441	1.530	-6.2	70	0.00
58		2,3,4,6-Tetrachlorophenol	0.268	0.239	10.8	56	0.00
59		Diethylphthalate	1.433	1.652	-15.3	75	0.00
60		4-Chlorophenyl-phenylether	0.590	0.644	-9.2	72	0.00
61		4-Nitroaniline	0.328	0.338	-3.0	63	0.00
62		Azobenzene	1.494	1.687	-12.9	74	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	71	0.00
64		4,6-Dinitro-2-methylphenol	0.104	0.105	-1.0	56	0.00
65	c	n-Nitrosodiphenylamine	0.713	0.774	-8.6	71	0.00
66		4-Bromophenyl-phenylether	0.203	0.222	-9.4	70	0.00
67		Hexachlorobenzene	0.214	0.220	-2.8	70	0.00
68		Atrazine	0.216	0.240	-11.1	69	0.00
69	C	Pentachlorophenol	0.082	0.068	17.1	53	0.00
70		Phenanthrene	1.247	1.296	-3.9	70	0.00
71		Anthracene	1.216	1.304	-7.2	73	0.00
72		Carbazole	1.180	1.277	-8.2	72	0.00
73		Di-n-butylphthalate	1.365	1.619	-18.6	77	0.00
74	C	Fluoranthene	1.278	1.376	-7.7	70	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	70	0.00
76		Benzidine	0.640	0.782	-22.2	86	0.00
77		Pyrene	1.371	1.487	-8.5	70	0.00
78	S	Terphenyl-d14	0.869	0.959	-10.4	72	0.00
79		Butylbenzylphthalate	0.621	0.755	-21.6	75	0.00
80		Benzo(a)anthracene	1.255	1.341	-6.9	69	0.00
81		3,3'-Dichlorobenzidine	0.399	0.435	-9.0	71	0.00
82		Chrysene	1.144	1.249	-9.2	71	0.00
83		Bis(2-ethylhexyl)phthalate	0.859	1.097	-27.7#	83	0.00
84	c	Di-n-octyl phthalate	1.491	1.919	-28.7#	82	0.00
85		Indeno(1,2,3-cd)pyrene	1.213	1.366	-12.6	72	0.00
86	I	Perylene-d12	1.000	1.000	0.0	69	0.00
87		Benzo(b)fluoranthene	1.347	1.557	-15.6	83	0.00

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88	Benzo(k)fluoranthene	1.177	1.224	-4.0	63	0.00
89 C	Benzo(a)pyrene	1.188	1.296	-9.1	71	0.00
90	Dibenzo(a,h)anthracene	1.090	1.228	-12.7	73	0.00
91	Benzo(g,h,i)perylene	1.084	1.227	-13.2	73	0.00

(#) = Out of Range SPCC's out = 0 CCC's out = 1