Evaluate Continuing Calibration Report

Data Path : V:\HPCHEM1\BNA F\DATA\BF112115\
Data File : BF083117.D
Acq On : 21 Nov 2015 18:05
Operator : UM/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF112115

Quant Time: Nov 22 09:26:55 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF112115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sun Nov 22 09:22:46 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	106	0.00
2		1,4-Dioxane	0.655	0.593	9.5	100	-0.02
3		Pyridine	1.731	1.793	-3.6	112	-0.11
4		n-Nitrosodimethylamine	0.798	0.841	-5.4	102	-0.03
5	S	2-Fluorophenol	1.323	1.321	0.2	108	-0.01
6		Aniline	2.092	2.145	-2.5	105	-0.01
7	S	Phenol-d6	1.712	1.696	0.9	110	-0.01
8		2-Chlorophenol	1.429	1.442	-0.9	110	0.00
9		Benzaldehyde	0.894	0.876	2.0	107	-0.01
10	C	Phenol	2.071	1.954	5.6	112	0.00
11		bis(2-Chloroethyl)ether	1.479	1.580	-6.8	116	0.00
12		1,3-Dichlorobenzene	1.632	1.586	2.8	106	0.00
13	C	1,4-Dichlorobenzene	1.651	1.650	0.1	110	0.00
14		1,2-Dichlorobenzene	1.500	1.424	5.1	100	0.00
15		Benzyl Alcohol	1.430	1.459	-2.0	107	-0.01
16		2,2'-oxybis(1-Chloropropane	2.306	2.394	-3.8	114	0.00
17		2-Methylphenol	1.183	1.152	2.6	107	-0.01
18		Hexachloroethane	0.581	0.568	2.2	104	-0.01
19	P	n-Nitroso-di-n-propylamine	1.191	1.237	-3.9	117	0.00
20		3+4-Methylphenols	1.495	1.540	-3.0	109	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	107	-0.01
22		Acetophenone	0.574	0.576	-0.3	114	0.00
23	S	Nitrobenzene-d5	0.438	0.423	3.4	105	0.00
24		Nitrobenzene	0.468	0.498	-6.4	117	0.00
25		Isophorone	0.829	0.846	-2.1	112	-0.01
26	C	2-Nitrophenol	0.206	0.191	7.3	93	0.00
27		2,4-Dimethylphenol	0.330	0.349	-5.8	119	0.00
28		bis(2-Chloroethoxy)methane	0.482	0.465	3.5	103	-0.01
29	C	2,4-Dichlorophenol	0.313	0.311	0.6	106	0.00
30		1,2,4-Trichlorobenzene	0.344	0.334	2.9	106	0.00
31		Naphthalene	1.079	1.012	6.2	103	-0.01
32		Benzoic acid	0.256	0.281	-9.8	109	-0.01
33		4-Chloroaniline	0.419	0.433	-3.3	114	0.00
34	C	Hexachlorobutadiene	0.203	0.203	0.0	110	0.00
35		Caprolactam	0.100	0.102	-2.0	112	-0.02
36	C	4-Chloro-3-methylphenol	0.363	0.366	-0.8	112	-0.01
37		2-Methylnaphthalene	0.701	0.719	-2.6	115	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	109	0.00
39		1,2,4,5-Tetrachlorobenzene	0.648	0.597	7.9	104	-0.01
40	P	Hexachlorocyclopentadiene	0.369	0.359	2.7	104	0.00
41	S	2,4,6-Tribromophenol	0.205	0.200	2.4	104	0.00
42	C	2,4,6-Trichlorophenol	0.464	0.437	5.8	104	-0.01
43		2,4,5-Trichlorophenol	0.475	0.485	-2.1	115	-0.01
44	S	2-Fluorobiphenyl	1.434	1.268	11.6	110	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.702	1.674	1.6	104	0.00
46		2-Chloronaphthalene	1.337	1.235	7.6	104	-0.01
47		2-Nitroaniline	0.473	0.507	-7.2	123	0.00
48		Acenaphthylene	2.072	1.987	4.1	106	0.00
49		Dimethylphthalate	1.540	1.522	1.2	110	0.00
50		2,6-Dinitrotoluene	0.345	0.337	2.3	118	0.00
51	C	Acenaphthene	1.261	1.191	5.6	104	-0.01
52		3-Nitroaniline	0.368	0.346	6.0	100	-0.01
53	P	2,4-Dinitrophenol	0.199	0.204	-2.5	110	0.00
54		Dibenzofuran	1.783	1.691	5.2	104	0.00
55	P	4-Nitrophenol	0.282	0.262	7.1	104	0.00
56		2,4-Dinitrotoluene	0.438	0.441	-0.7	118	0.00
57		Fluorene	1.474	1.313	10.9	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.389	0.384	1.3	106	0.00
59		Diethylphthalate	1.533	1.446	5.7	104	-0.01
60		4-Chlorophenyl-phenylether	0.675	0.653	3.3	112	0.00
61		4-Nitroaniline	0.363	0.364	-0.3	104	-0.01
62		Azobenzene	1.749	1.771	-1.3	115	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	105	0.00
64		4,6-Dinitro-2-methylphenol	0.138	0.154	-11.6	108	0.00
65	c	n-Nitrosodiphenylamine	0.680	0.704	-3.5	115	0.00
66		4-Bromophenyl-phenylether	0.222	0.217	2.3	102	0.00
67		Hexachlorobenzene	0.244	0.251	-2.9	110	0.00
68		Atrazine	0.198	0.184	7.1	97	-0.01
69	C	Pentachlorophenol	0.159	0.160	-0.6	103	0.00
70		Phenanthrene	1.152	1.166	-1.2	108	0.00
71		Anthracene	1.176	1.081	8.1	104	0.00
72		Carbazole	1.045	1.021	2.3	102	0.00
73		Di-n-butylphthalate	1.275	1.275	0.0	108	0.00
74	C	Fluoranthene	1.179	1.136	3.6	110	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
76		Benzidine	0.524	0.632	-20.6	112	0.00
77		Pyrene	1.366	1.517	-11.1	116	0.00
78	S	Terphenyl-d14	0.849	0.805	5.2	96	-0.01
79		Butylbenzylphthalate	0.647	0.696	-7.6	108	0.00
80		Benzo(a)anthracene	1.237	1.304	-5.4	107	0.00
81		3,3'-Dichlorobenzidine	0.431	0.425	1.4	101	-0.01
82		Chrysene	1.147	1.287	-12.2	119	0.00
83		Bis(2-ethylhexyl)phthalate	0.845	0.941	-11.4	116	0.00
84	c	Di-n-octyl phthalate	1.454	1.539	-5.8	110	-0.01
85		Indeno(1,2,3-cd)pyrene	1.043	1.080	-3.5	114	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	112	0.00
87		Benzo(b)fluoranthene	1.367	1.131	17.3	93	-0.01

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.018	1.186	-16.5	142	0.00
89 C	Benzo(a)pyrene	1.128	1.052	6.7	108	-0.01
90	Dibenzo(a,h)anthracene	1.025	0.981	4.3	114	-0.01
91	Benzo(g,h,i)perylene	1.000	0.928	7.2	110	-0.01

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