Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF072315\
Data File : BF080597.D
Acq On : 23 Jul 2015 17:47
Operator : TP/UM
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF072315

Quant Time: Jul 24 06:56:31 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF072315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 24 01:01:51 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	104	0.00
2		1,4-Dioxane	0.552	0.548	0.7	97	0.00
3		Pyridine	1.466	1.546	-5.5	100	0.00
4		n-Nitrosodimethylamine	0.927	0.958	-3.3	99	0.00
5	S	2-Fluorophenol	1.158	1.194	-3.1	100	0.00
6		Aniline	2.069	2.216	-7.1	102	0.00
7	S	Phenol-d6	1.416	1.567	-10.7	102	0.00
8		2-Chlorophenol	1.244	1.277	-2.7	99	0.00
9		Benzaldehyde	1.047	1.042	0.5	97	0.00
10	C	Phenol	1.694	1.881	-11.0	103	0.00
11		bis(2-Chloroethyl)ether	1.176	1.204	-2.4	104	0.00
12		1,3-Dichlorobenzene	1.577	1.620	-2.7	103	0.00
13	C	1,4-Dichlorobenzene	1.548	1.611	-4.1	104	0.00
14		1,2-Dichlorobenzene	1.435	1.470	-2.4	100	0.00
15		Benzyl Alcohol	1.247	1.289	-3.4	102	0.00
16		2,2'-oxybis(1-Chloropropane	2.384	2.427	-1.8	99	0.00
17		2-Methylphenol	0.980	0.996	-1.6	99	0.00
18		Hexachloroethane	0.535	0.555	-3.7	102	0.00
19	P	n-Nitroso-di-n-propylamine	0.959	1.011	-5.4	106	0.00
20		3+4-Methylphenols	1.355	1.414	-4.4	104	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
22		Acetophenone	0.487	0.484	0.6	96	0.00
23	S	Nitrobenzene-d5	0.353	0.378	-7.1	102	0.00
24		Nitrobenzene	0.377	0.404	-7.2	102	0.00
25		Isophorone	0.654	0.678	-3.7	100	0.00
26	C	2-Nitrophenol	0.135	0.147	-8.9	110	0.00
27		2,4-Dimethylphenol	0.298	0.301	-1.0	98	0.00
28		bis(2-Chloroethoxy)methane	0.361	0.351	2.8	99	0.00
29	C	2,4-Dichlorophenol	0.244	0.250	-2.5	104	0.00
30		1,2,4-Trichlorobenzene	0.317	0.310	2.2	101	0.00
31		Naphthalene	1.000	0.980	2.0	99	0.00
32		Benzoic acid	0.127	0.141	-11.0	115	0.00
33		4-Chloroaniline	0.392	0.387	1.3	98	0.00
34	C	Hexachlorobutadiene	0.185	0.191	-3.2	105	0.00
35		Caprolactam	0.066	0.070	-6.1	97	0.00
36	C	4-Chloro-3-methylphenol	0.290	0.311	-7.2	99	0.00
37		2-Methylnaphthalene	0.560	0.561	-0.2	99	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	101	0.00
39		1,2,4,5-Tetrachlorobenzene	0.676	0.707	-4.6	103	0.00
40	P	Hexachlorocyclopentadiene	0.356	0.388	-9.0	108	0.00
41	S	2,4,6-Tribromophenol	0.158	0.159	-0.6	97	0.00
42	C	2,4,6-Trichlorophenol	0.379	0.384	-1.3	101	0.00
43		2,4,5-Trichlorophenol	0.383	0.415	-8.4	107	0.00
44	S	2-Fluorobiphenyl	1.378	1.411	-2.4	104	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.628	1.640	-0.7	101	0.00
46		2-Chloronaphthalene	1.317	1.344	-2.1	102	0.00
47		2-Nitroaniline	0.402	0.409	-1.7	102	0.00
48		Acenaphthylene	1.951	1.959	-0.4	101	0.00
49		Dimethylphthalate	1.471	1.537	-4.5	100	0.00
50		2,6-Dinitrotoluene	0.275	0.299	-8.7	103	0.00
51	C	Acenaphthene	1.172	1.174	-0.2	101	0.00
52		3-Nitroaniline	0.299	0.296	1.0	102	0.00
53	P	2,4-Dinitrophenol	0.076	0.074	2.6	103	0.00
54		Dibenzofuran	1.694	1.641	3.1	96	0.00
55	P	4-Nitrophenol	0.216	0.248	-14.8	118	0.00
56		2,4-Dinitrotoluene	0.345	0.376	-9.0	102	0.00
57		Fluorene	1.360	1.341	1.4	95	0.00
58		2,3,4,6-Tetrachlorophenol	0.298	0.298	0.0	92	0.00
59		Diethylphthalate	1.373	1.409	-2.6	111	0.00
60		4-Chlorophenyl-phenylether	0.611	0.607	0.7	101	0.00
61		4-Nitroaniline	0.301	0.291	3.3	98	0.00
62		Azobenzene	1.439	1.508	-4.8	100	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	96	0.00
64		4,6-Dinitro-2-methylphenol	0.065	0.068	-4.6	107	0.00
65	c	n-Nitrosodiphenylamine	0.656	0.719	-9.6	104	0.00
66		4-Bromophenyl-phenylether	0.203	0.217	-6.9	102	0.00
67		Hexachlorobenzene	0.230	0.222	3.5	95	0.00
68		Atrazine	0.180	0.194	-7.8	95	0.00
69	C	Pentachlorophenol	0.107	0.118	-10.3	106	0.00
70		Phenanthrene	1.129	1.198	-6.1	100	0.00
71		Anthracene	1.057	1.101	-4.2	101	0.00
72		Carbazole	0.914	0.927	-1.4	95	0.00
73		Di-n-butylphthalate	1.062	1.067	-0.5	91	-0.01
74	C	Fluoranthene	1.010	1.064	-5.3	98	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	104	-0.07
76		Benzidine	0.588	0.551	6.3	94	-0.02
77		Pyrene	1.342	1.285	4.2	99	-0.02
78	S	Terphenyl-d14	0.956	0.931	2.6	101	-0.02
79		Butylbenzylphthalate	0.444	0.428	3.6	94	-0.04
80		Benzo(a)anthracene	1.128	1.131	-0.3	102	-0.07
81		3,3'-Dichlorobenzidine	0.322	0.315	2.2	93	-0.07
82		Chrysene	1.122	1.069	4.7	98	-0.08
83		Bis(2-ethylhexyl)phthalate	0.656	0.648	1.2	99	-0.08
84	c	Di-n-octyl phthalate	0.829	0.799	3.6	92	-0.11
85		Indeno(1,2,3-cd)pyrene	0.849	0.797	6.1	91	-0.14
86	I	Perylene-d12	1.000	1.000	0.0	101	-0.14
87		Benzo(b)fluoranthene	1.216	1.176	3.3	94	-0.13

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88	Benzo(k)fluoranthene	1.211	1.320	-9.0	101	-0.14
89 C	Benzo(a)pyrene	1.079	1.141	-5.7	100	-0.13
90	Dibenzo(a,h)anthracene	1.019	0.998	2.1	91	-0.15
91	Benzo(g,h,i)perylene	1.041	1.015	2.5	91	-0.15

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