Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF071715\
Data File : BF080543.D
Acq On : 17 Jul 2015 19:57
Operator : TP/UM
Sample : SSTDICV040
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF071715

Quant Time: Jul 20 14:11:02 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF071715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 20 13:29:49 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	96	0.00
2		1,4-Dioxane	0.537	0.607	-13.0	108	0.00
3		Pyridine	1.097	1.105	-0.7	101	0.00
4		n-Nitrosodimethylamine	0.707	0.831	-17.5	105	0.00
5	S	2-Fluorophenol	1.283	1.284	-0.1	95	0.00
6		Aniline	1.924	1.956	-1.7	98	0.00
7	S	Phenol-d6	1.460	1.531	-4.9	99	0.00
8		2-Chlorophenol	1.338	1.356	-1.3	100	0.00
9		Benzaldehyde	0.824	0.932	-13.1	111	0.00
10	C	Phenol	1.688	1.655	2.0	98	-0.01
11		bis(2-Chloroethyl)ether	1.276	1.199	6.0	101	0.00
12		1,3-Dichlorobenzene	1.507	1.579	-4.8	100	0.00
13	C	1,4-Dichlorobenzene	1.659	1.693	-2.0	99	0.00
14		1,2-Dichlorobenzene	1.526	1.505	1.4	98	0.00
15		Benzyl Alcohol	0.932	0.900	3.4	98	0.00
16		2,2'-oxybis(1-Chloropropane	2.171	2.027	6.6	96	0.00
17		2-Methylphenol	1.021	0.983	3.7	100	0.00
18		Hexachloroethane	0.532	0.496	6.8	97	0.00
19	P	n-Nitroso-di-n-propylamine	0.869	0.801	7.8	95	0.00
20		3+4-Methylphenols	1.474	1.464	0.7	97	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
22		Acetophenone	0.429	0.426	0.7	95	0.00
23	S	Nitrobenzene-d5	0.314	0.351	-11.8	101	0.00
24		Nitrobenzene	0.354	0.336	5.1	98	0.00
25		Isophorone	0.599	0.613	-2.3	102	0.00
26	C	2-Nitrophenol	0.174	0.176	-1.1	104	0.00
27		2,4-Dimethylphenol	0.298	0.329	-10.4	101	0.00
28		bis(2-Chloroethoxy)methane	0.332	0.318	4.2	98	0.00
29	C	2,4-Dichlorophenol	0.252	0.268	-6.3	100	0.00
30		1,2,4-Trichlorobenzene	0.322	0.324	-0.6	97	0.00
31		Naphthalene	0.990	1.044	-5.5	97	0.00
32		Benzoic acid	0.133	0.128	3.8	99	0.00
33		4-Chloroaniline	0.394	0.441	-11.9	102	0.00
34	C	Hexachlorobutadiene	0.150	0.155	-3.3	97	0.00
35		Caprolactam	0.061	0.061	0.0	102	0.00
36	C	4-Chloro-3-methylphenol	0.260	0.272	-4.6	96	0.00
37		2-Methylnaphthalene	0.608	0.635	-4.4	100	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	90	0.00
39		1,2,4,5-Tetrachlorobenzene	0.632	0.669	-5.9	99	0.00
40	P	Hexachlorocyclopentadiene	0.298	0.338	-13.4	105	0.00
41	S	2,4,6-Tribromophenol	0.179	0.189	-5.6	93	0.00
42	C	2,4,6-Trichlorophenol	0.364	0.418	-14.8	97	0.00
43		2,4,5-Trichlorophenol	0.398	0.422	-6.0	99	0.00
44	S	2-Fluorobiphenyl	1.424	1.515	-6.4	98	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.658	1.737	-4.8	100	0.00
46		2-Chloronaphthalene	1.222	1.343	-9.9	99	0.00
47		2-Nitroaniline	0.293	0.301	-2.7	92	0.00
48		Acenaphthylene	2.023	2.129	-5.2	104	0.00
49		Dimethylphthalate	1.331	1.360	-2.2	99	0.00
50		2,6-Dinitrotoluene	0.259	0.270	-4.2	102	0.01
51	C	Acenaphthene	1.236	1.236	0.0	96	0.00
52		3-Nitroaniline	0.294	0.327	-11.2	102	0.00
53	P	2,4-Dinitrophenol	0.113	0.118	-4.4	102	0.00
54		Dibenzofuran	1.721	1.750	-1.7	99	0.00
55	P	4-Nitrophenol	0.218	0.230	-5.5	101	0.00
56		2,4-Dinitrotoluene	0.353	0.426	-20.7	108	0.00
57		Fluorene	1.364	1.383	-1.4	100	0.00
58		2,3,4,6-Tetrachlorophenol	0.271	0.291	-7.4	99	0.00
59		Diethylphthalate	1.256	1.305	-3.9	99	0.00
60		4-Chlorophenyl-phenylether	0.611	0.609	0.3	99	0.00
61		4-Nitroaniline	0.289	0.350	-21.1	103	0.00
62		Azobenzene	1.264	1.267	-0.2	97	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	97	0.00
64		4,6-Dinitro-2-methylphenol	0.097	0.088	9.3	100	0.00
65	c	n-Nitrosodiphenylamine	0.659	0.648	1.7	94	0.00
66		4-Bromophenyl-phenylether	0.183	0.176	3.8	100	0.00
67		Hexachlorobenzene	0.222	0.227	-2.3	101	0.00
68		Atrazine	0.156	0.158	-1.3	106	0.01
69	C	Pentachlorophenol	0.090	0.086	4.4	100	0.00
70		Phenanthrene	1.090	1.075	1.4	98	0.00
71		Anthracene	1.058	1.045	1.2	104	0.00
72		Carbazole	0.948	0.910	4.0	100	0.00
73		Di-n-butylphthalate	0.980	1.111	-13.4	115	0.00
74	C	Fluoranthene	1.019	0.994	2.5	96	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	95	0.06
76		Benzidine	0.535	0.605	-13.1	112	0.02
77		Pyrene	1.291	1.251	3.1	97	0.02
78	S	Terphenyl-d14	0.896	0.916	-2.2	97	0.02
79		Butylbenzylphthalate	0.402	0.448	-11.4	102	0.03
80		Benzo(a)anthracene	1.113	1.145	-2.9	100	0.06
81		3,3'-Dichlorobenzidine	0.352	0.379	-7.7	103	0.06
82		Chrysene	1.115	1.144	-2.6	101	0.06
83		Bis(2-ethylhexyl)phthalate	0.615	0.666	-8.3	98	0.07
84	c	Di-n-octyl phthalate	0.873	0.961	-10.1	103	0.09
85		Indeno(1,2,3-cd)pyrene	1.149	1.077	6.3	98	0.14
86	I	Perylene-d12	1.000	1.000	0.0	103	0.11
87		Benzo(b)fluoranthene	1.240	1.168	5.8	102	0.11

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88	Benzo(k)fluoranthene	1.073	1.100	-2.5	94	0.14
89 C	Benzo(a)pyrene	1.056	1.096	-3.8	106	0.11
90	Dibenzo(a,h)anthracene	1.040	1.003	3.6	101	0.14
91	Benzo(g,h,i)perylene	1.050	0.982	6.5	98	0.13

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