Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF102615\
Data File : BF082525.D
Acq On : 26 Oct 2015 21:49
Operator : UM/IZ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF102615

Quant Time: Oct 27 06:17:36 2015
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF102615.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Oct 27 02:43:58 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	89	0.00
2		1,4-Dioxane	0.508	0.496	2.4	94	0.00
3		Pyridine	1.185	1.249	-5.4	90	-0.01
4		n-Nitrosodimethylamine	0.521	0.540	-3.6	91	-0.01
5	S	2-Fluorophenol	1.013	1.038	-2.5	92	0.00
6		Aniline	1.664	1.664	0.0	91	0.00
7	S	Phenol-d6	1.320	1.281	3.0	91	0.00
8		2-Chlorophenol	1.188	1.240	-4.4	94	0.00
9		Benzaldehyde	0.688	0.690	-0.3	91	0.00
10	C	Phenol	1.499	1.364	9.0	90	-0.01
11		bis(2-Chloroethyl)ether	1.190	1.149	3.4	93	0.00
12		1,3-Dichlorobenzene	1.417	1.403	1.0	93	0.00
13	C	1,4-Dichlorobenzene	1.458	1.457	0.1	92	0.00
14		1,2-Dichlorobenzene	1.354	1.257	7.2	93	0.00
15		Benzyl Alcohol	0.919	0.889	3.3	89	0.00
16		2,2'-oxybis(1-Chloropropane	1.611	1.624	-0.8	92	0.00
17		2-Methylphenol	0.953	0.978	-2.6	90	0.00
18		Hexachloroethane	0.496	0.504	-1.6	92	0.00
19	P	n-Nitroso-di-n-propylamine	0.860	0.850	1.2	95	0.00
20		3+4-Methylphenols	1.218	1.234	-1.3	93	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
22		Acetophenone	0.434	0.428	1.4	92	0.00
23	S	Nitrobenzene-d5	0.310	0.296	4.5	91	0.00
24		Nitrobenzene	0.346	0.354	-2.3	91	0.00
25		Isophorone	0.616	0.585	5.0	90	0.00
26	C	2-Nitrophenol	0.167	0.160	4.2	88	-0.01
27		2,4-Dimethylphenol	0.272	0.260	4.4	91	0.00
28		bis(2-Chloroethoxy)methane	0.361	0.340	5.8	91	0.00
29	C	2,4-Dichlorophenol	0.271	0.272	-0.4	92	0.00
30		1,2,4-Trichlorobenzene	0.307	0.296	3.6	92	0.00
31		Naphthalene	0.914	0.844	7.7	91	-0.01
32		Benzoic acid	0.152	0.124	18.4	72	0.00
33		4-Chloroaniline	0.354	0.362	-2.3	90	0.00
34	C	Hexachlorobutadiene	0.177	0.167	5.6	92	0.00
35		Caprolactam	0.075	0.079	-5.3	101	0.00
36	C	4-Chloro-3-methylphenol	0.264	0.275	-4.2	92	0.00
37		2-Methylnaphthalene	0.617	0.607	1.6	92	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	93	0.00
39		1,2,4,5-Tetrachlorobenzene	0.604	0.579	4.1	90	0.00
40	P	Hexachlorocyclopentadiene	0.115	0.122	-6.1	89	0.00
41	S	2,4,6-Tribromophenol	0.150	0.152	-1.3	93	0.00
42	C	2,4,6-Trichlorophenol	0.376	0.378	-0.5	91	0.00
43		2,4,5-Trichlorophenol	0.395	0.408	-3.3	91	0.00
44	S	2-Fluorobiphenyl	1.258	1.215	3.4	96	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.488	1.548	-4.0	90	0.00
46		2-Chloronaphthalene	1.173	1.209	-3.1	94	0.00
47		2-Nitroaniline	0.341	0.323	5.3	90	0.00
48		Acenaphthylene	1.879	1.800	4.2	93	0.00
49		Dimethylphthalate	1.392	1.409	-1.2	92	0.00
50		2,6-Dinitrotoluene	0.305	0.317	-3.9	90	0.00
51	C	Acenaphthene	1.078	1.042	3.3	93	0.00
52		3-Nitroaniline	0.331	0.348	-5.1	90	0.00
53	P	2,4-Dinitrophenol	0.112	0.108	3.6	88	0.00
54		Dibenzofuran	1.572	1.477	6.0	91	0.00
55	P	4-Nitrophenol	0.113	0.120	-6.2	88	0.00
56		2,4-Dinitrotoluene	0.370	0.366	1.1	94	0.00
57		Fluorene	1.291	1.225	5.1	94	0.00
58		2,3,4,6-Tetrachlorophenol	0.302	0.326	-7.9	99	0.00
59		Diethylphthalate	1.354	1.333	1.6	90	0.00
60		4-Chlorophenyl-phenylether	0.607	0.556	8.4	90	-0.01
61		4-Nitroaniline	0.332	0.353	-6.3	90	0.00
62		Azobenzene	1.308	1.191	8.9	90	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	92	0.00
64		4,6-Dinitro-2-methylphenol	0.117	0.121	-3.4	94	0.00
65	c	n-Nitrosodiphenylamine	0.605	0.546	9.8	90	0.00
66		4-Bromophenyl-phenylether	0.188	0.173	8.0	90	0.00
67		Hexachlorobenzene	0.200	0.208	-4.0	95	0.00
68		Atrazine	0.187	0.194	-3.7	95	0.00
69	C	Pentachlorophenol	0.068	0.069	-1.5	85	0.00
70		Phenanthrene	1.029	0.930	9.6	94	0.00
71		Anthracene	1.037	1.025	1.2	90	0.00
72		Carbazole	0.914	0.916	-0.2	93	0.00
73		Di-n-butylphthalate	1.090	0.990	9.2	92	0.00
74	C	Fluoranthene	1.057	1.000	5.4	94	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	87	-0.01
76		Benzidine	0.549	0.590	-7.5	92	0.00
77		Pyrene	1.230	1.204	2.1	90	0.00
78	S	Terphenyl-d14	0.694	0.738	-6.3	93	0.00
79		Butylbenzylphthalate	0.517	0.510	1.4	91	0.00
80		Benzo(a)anthracene	1.072	1.089	-1.6	94	-0.01
81		3,3'-Dichlorobenzidine	0.372	0.389	-4.6	88	0.00
82		Chrysene	1.013	1.127	-11.3	101	0.00
83		Bis(2-ethylhexyl)phthalate	0.693	0.706	-1.9	89	-0.01
84	c	Di-n-octyl phthalate	1.143	1.263	-10.5	98	-0.01
85		Indeno(1,2,3-cd)pyrene	0.931	0.959	-3.0	95	-0.02
86	I	Perylene-d12	1.000	1.000	0.0	95	-0.01
87		Benzo(b)fluoranthene	1.188	1.061	10.7	88	-0.02

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88	Benzo(k)fluoranthene	0.968	1.038	-7.2	105	-0.02
89 C	Benzo(a)pyrene	1.029	0.996	3.2	94	-0.02
90	Dibenzo(a,h)anthracene	0.861	0.843	2.1	96	-0.02
91	Benzo(g,h,i)perylene	0.852	0.789	7.4	93	-0.02

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