Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF040216\
Data File : BF085970.D
Acq On : 1 Apr 2016 19:18
Operator : UM/SJ
Sample : SSTDICV040
Misc :
ALS Vial : 10 Sample Multiplier: 1

Instrument : BNA_F
ClientSampleId : ICVBF040216

Quant Time: Apr 01 19:46:32 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF040216.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Apr 01 19:44:06 2016
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	98	0.00
2		1,4-Dioxane	0.555	0.540	2.7	100	0.00
3		Pyridine	1.242	1.456	-17.2	104	0.00
4		n-Nitrosodimethylamine	0.546	0.570	-4.4	101	0.00
5	S	2-Fluorophenol	1.170	1.218	-4.1	101	0.00
6		Aniline	1.950	2.012	-3.2	100	0.00
7	S	Phenol-d6	1.486	1.563	-5.2	101	0.00
8		2-Chlorophenol	1.395	1.402	-0.5	101	0.00
9		Benzaldehyde	0.800	0.804	-0.5	98	0.00
10	C	Phenol	1.695	1.861	-9.8	101	0.00
11		bis(2-Chloroethyl)ether	1.342	1.228	8.5	100	0.00
12		1,3-Dichlorobenzene	1.479	1.505	-1.8	103	0.00
13	C	1,4-Dichlorobenzene	1.524	1.517	0.5	102	0.00
14		1,2-Dichlorobenzene	1.402	1.467	-4.6	101	0.00
15		Benzyl Alcohol	0.962	0.951	1.1	101	0.00
16		2,2'-oxybis(1-Chloropropane	1.640	1.649	-0.5	100	0.00
17		2-Methylphenol	1.100	1.088	1.1	103	0.00
18		Hexachloroethane	0.560	0.563	-0.5	102	0.00
19	P	n-Nitroso-di-n-propylamine	0.921	0.861	6.5	100	0.00
20		3+4-Methylphenols	1.338	1.321	1.3	100	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	97	0.00
22		Acetophenone	0.470	0.480	-2.1	104	0.00
23	S	Nitrobenzene-d5	0.339	0.332	2.1	102	0.00
24		Nitrobenzene	0.371	0.389	-4.9	103	0.00
25		Isophorone	0.669	0.671	-0.3	100	0.00
26	C	2-Nitrophenol	0.178	0.178	0.0	101	0.00
27		2,4-Dimethylphenol	0.319	0.307	3.8	103	0.00
28		bis(2-Chloroethoxy)methane	0.393	0.372	5.3	101	0.00
29	C	2,4-Dichlorophenol	0.270	0.269	0.4	100	0.00
30		1,2,4-Trichlorobenzene	0.303	0.301	0.7	101	0.00
31		Naphthalene	1.006	0.990	1.6	100	0.00
32		Benzoic acid	0.234	0.249	-6.4	101	0.00
33		4-Chloroaniline	0.406	0.396	2.5	100	0.00
34	C	Hexachlorobutadiene	0.163	0.163	0.0	100	0.00
35		Caprolactam	0.086	0.098	-14.0	104	0.01
36	C	4-Chloro-3-methylphenol	0.303	0.313	-3.3	100	0.00
37		2-Methylnaphthalene	0.657	0.632	3.8	101	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	99	0.00
39		1,2,4,5-Tetrachlorobenzene	0.601	0.595	1.0	99	0.00
40	P	Hexachlorocyclopentadiene	0.163	0.180	-10.4	107	0.00
41	S	2,4,6-Tribromophenol	0.188	0.194	-3.2	101	0.00
42	C	2,4,6-Trichlorophenol	0.386	0.386	0.0	99	0.00
43		2,4,5-Trichlorophenol	0.392	0.398	-1.5	103	0.00
44	S	2-Fluorobiphenyl	1.203	1.275	-6.0	102	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.724	1.736	-0.7	102	0.00
46		2-Chloronaphthalene	1.250	1.331	-6.5	104	0.00
47		2-Nitroaniline	0.366	0.389	-6.3	102	0.00
48		Acenaphthylene	2.002	2.111	-5.4	101	0.00
49		Dimethylphthalate	1.482	1.442	2.7	104	0.00
50		2,6-Dinitrotoluene	0.314	0.340	-8.3	100	0.00
51	C	Acenaphthene	1.191	1.213	-1.8	99	0.00
52		3-Nitroaniline	0.378	0.411	-8.7	100	0.00
53	P	2,4-Dinitrophenol	0.127	0.160	-26.0#	102	0.00
54		Dibenzofuran	1.573	1.603	-1.9	100	0.00
55	P	4-Nitrophenol	0.223	0.227	-1.8	102	0.00
56		2,4-Dinitrotoluene	0.396	0.392	1.0	104	0.00
57		Fluorene	1.344	1.382	-2.8	98	0.00
58		2,3,4,6-Tetrachlorophenol	0.292	0.288	1.4	99	0.00
59		Diethylphthalate	1.496	1.399	6.5	100	0.00
60		4-Chlorophenyl-phenylether	0.626	0.616	1.6	98	0.00
61		4-Nitroaniline	0.372	0.386	-3.8	103	0.01
62		Azobenzene	1.433	1.443	-0.7	101	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.00
64		4,6-Dinitro-2-methylphenol	0.121	0.126	-4.1	101	0.00
65	c	n-Nitrosodiphenylamine	0.625	0.668	-6.9	101	0.00
66		4-Bromophenyl-phenylether	0.204	0.220	-7.8	101	0.00
67		Hexachlorobenzene	0.211	0.231	-9.5	103	0.00
68		Atrazine	0.202	0.197	2.5	101	0.00
69	C	Pentachlorophenol	0.099	0.106	-7.1	99	0.00
70		Phenanthrene	1.091	1.143	-4.8	100	0.00
71		Anthracene	1.143	1.077	5.8	101	0.00
72		Carbazole	0.982	0.964	1.8	101	0.00
73		Di-n-butylphthalate	1.217	1.306	-7.3	102	0.00
74	C	Fluoranthene	1.135	1.122	1.1	103	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	98	0.00
76		Benzidine	0.706	0.752	-6.5	103	0.00
77		Pyrene	1.409	1.359	3.5	99	0.00
78	S	Terphenyl-d14	0.851	0.839	1.4	104	0.00
79		Butylbenzylphthalate	0.635	0.661	-4.1	102	-0.01
80		Benzo(a)anthracene	1.179	1.169	0.8	101	0.00
81		3,3'-Dichlorobenzidine	0.861	0.843	2.1	91	-0.01
82		Chrysene	1.136	1.163	-2.4	97	0.00
83		Bis(2-ethylhexyl)phthalate	0.842	0.854	-1.4	100	0.00
84	c	Di-n-octyl phthalate	1.345	1.470	-9.3	100	0.00
85		Indeno(1,2,3-cd)pyrene	0.921	0.888	3.6	95	0.00
86	I	Perylene-d12	1.000	1.000	0.0	94	0.00
87		Benzo(b)fluoranthene	1.258	1.524	-21.1	101	-0.01

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88	Benzo(k)fluoranthene	1.114	0.913	18.0	92	0.00
89 C	Benzo(a)pyrene	1.115	1.131	-1.4	99	0.00
90	Dibenzo(a,h)anthracene	0.897	0.881	1.8	94	0.00
91	Benzo(g,h,i)perylene	0.868	0.849	2.2	95	-0.01

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