Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF082417\
Data File : BF098001.D
Acq On : 24 Aug 2017 14:50
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Aug 25 01:13:52 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF081517.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 21 15:59:04 2017
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	94	-0.01
2		1,4-Dioxane	0.733	0.735	-0.3	95	-0.02
3		Pyridine	2.027	2.031	-0.2	94	-0.02
4		n-Nitrosodimethylamine	0.780	0.760	2.6	88	-0.02
5	S	2-Fluorophenol	1.315	1.322	-0.5	95	-0.01
6		Aniline	2.510	2.439	2.8	92	-0.01
7	S	Phenol-d6	1.787	1.788	-0.1	95	-0.01
8		2-Chlorophenol	1.387	1.395	-0.6	94	0.00
9		Benzaldehyde	1.132	1.071	5.4	88	-0.01
10	C	Phenol	2.089	2.115	-1.2	92	0.00
11		bis(2-Chloroethyl)ether	1.577	1.533	2.8	92	-0.01
12		1,3-Dichlorobenzene	1.492	1.472	1.3	94	-0.01
13	C	1,4-Dichlorobenzene	1.517	1.489	1.8	93	-0.01
14		1,2-Dichlorobenzene	1.399	1.365	2.4	92	-0.01
15		Benzyl Alcohol	1.338	1.308	2.2	91	0.00
16		2,2'-oxybis(1-Chloropropane	2.122	1.978	6.8	88	-0.01
17		2-Methylphenol	1.222	1.202	1.6	92	-0.01
18		Hexachloroethane	0.595	0.599	-0.7	93	-0.01
19	P	n-Nitroso-di-n-propylamine	1.223	1.120	8.4	86	-0.01
20		3+4-Methylphenols	1.493	1.416	5.2	89	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	92	0.00
22		Acetophenone	0.528	0.496	6.1	89	0.00
23	S	Nitrobenzene-d5	0.368	0.412	-12.0	100	-0.01
24		Nitrobenzene	0.410	0.446	-8.8	94	-0.01
25		Isophorone	0.766	0.743	3.0	89	-0.01
26	C	2-Nitrophenol	0.137	0.176	-28.5#	113	-0.01
27		2,4-Dimethylphenol	0.319	0.321	-0.6	94	-0.01
28		bis(2-Chloroethoxy)methane	0.503	0.489	2.8	89	0.00
29	C	2,4-Dichlorophenol	0.265	0.273	-3.0	93	0.00
30		1,2,4-Trichlorobenzene	0.294	0.290	1.4	91	0.00
31		Naphthalene	1.038	1.020	1.7	92	0.00
32		Benzoic acid	0.212	0.261	-23.1	115	-0.02
33		4-Chloroaniline	0.438	0.433	1.1	92	0.00
34	C	Hexachlorobutadiene	0.172	0.169	1.7	91	-0.01
35		Caprolactam	0.090	0.098	-8.9	97	-0.02
36	C	4-Chloro-3-methylphenol	0.341	0.342	-0.3	90	0.00
37		2-Methylnaphthalene	0.637	0.630	1.1	92	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	90	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.614	0.614	0.0	92	0.00
40	P	Hexachlorocyclopentadiene	0.295	0.289	2.0	84	0.00
41	S	2,4,6-Tribromophenol	0.174	0.188	-8.0	95	-0.01
42	C	2,4,6-Trichlorophenol	0.412	0.429	-4.1	93	0.00
43		2,4,5-Trichlorophenol	0.409	0.432	-5.6	93	0.00
44	S	2-Fluorobiphenyl	1.361	1.293	5.0	90	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.824	1.789	1.9	90	0.00
46		2-Chloronaphthalene	1.348	1.330	1.3	92	0.00
47		2-Nitroaniline	0.452	0.541	-19.7	104	-0.01
48		Acenaphthylene	2.116	2.081	1.7	89	-0.01
49		Dimethylphthalate	1.462	1.451	0.8	91	-0.01
50		2,6-Dinitrotoluene	0.292	0.321	-9.9	96	0.00
51	C	Acenaphthene	1.335	1.265	5.2	87	-0.01
52		3-Nitroaniline	0.376	0.430	-14.4	102	-0.01
53	P	2,4-Dinitrophenol	0.106	0.160	-50.9#	134	0.00
54		Dibenzofuran	1.789	1.755	1.9	91	-0.01
55	P	4-Nitrophenol	0.300	0.330	-10.0	96	0.00
56		2,4-Dinitrotoluene	0.366	0.421	-15.0	99	0.00
57		Fluorene	1.383	1.371	0.9	93	0.00
58		2,3,4,6-Tetrachlorophenol	0.317	0.326	-2.8	92	-0.01
59		Diethylphthalate	1.529	1.479	3.3	87	0.00
60		4-Chlorophenyl-phenylether	0.642	0.629	2.0	90	-0.01
61		4-Nitroaniline	0.358	0.430	-20.1	105	-0.01
62		Azobenzene	1.816	1.738	4.3	87	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	94	0.00
64		4,6-Dinitro-2-methylphenol	0.083	0.113	-36.1#	127	0.00
65	c	n-Nitrosodiphenylamine	0.681	0.656	3.7	92	0.00
66		4-Bromophenyl-phenylether	0.208	0.204	1.9	93	-0.01
67		Hexachlorobenzene	0.212	0.204	3.8	92	-0.01
68		Atrazine	0.181	0.169	6.6	89	-0.01
69	C	Pentachlorophenol	0.123	0.125	-1.6	90	0.00
70		Phenanthrene	1.066	1.014	4.9	92	-0.01
71		Anthracene	1.081	1.046	3.2	93	-0.01
72		Carbazole	1.026	1.013	1.3	96	0.00
73		Di-n-butylphthalate	1.182	1.193	-0.9	94	0.00
74	C	Fluoranthene	1.112	1.114	-0.2	96	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
76		Benzidine	0.656	0.521	20.6	86	-0.01
77		Pyrene	1.636	1.549	5.3	97	-0.01
78	S	Terphenyl-d14	0.959	0.885	7.7	96	0.00
79		Butylbenzylphthalate	0.627	0.661	-5.4	105	-0.01
80		Benzo(a)anthracene	1.199	1.119	6.7	97	0.00
81		3,3'-Dichlorobenzidine	0.404	0.410	-1.5	99	0.00
82		Chrysene	1.192	1.126	5.5	99	-0.01
83		Bis(2-ethylhexyl)phthalate	0.695	0.751	-8.1	103	0.00
84	c	Di-n-octyl phthalate	1.209	1.299	-7.4	104	-0.01
85		Indeno(1,2,3-cd)pyrene	0.877	0.779	11.2	94	0.00
86	I	Perylene-d12	1.000	1.000	0.0	92	0.00
87		Benzo(b)fluoranthene	1.261	1.330	-5.5	96	0.00

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88	Benzo(k)fluoranthene	1.184	1.205	-1.8	91	-0.01
89 C	Benzo(a)pyrene	1.116	1.147	-2.8	92	-0.01
90	Dibenzo(a,h)anthracene	0.909	0.968	-6.5	95	-0.01
91	Benzo(g,h,i)perylene	0.948	1.031	-8.8	98	-0.02

(#) = Out of Range SPCC's out = 0 CCC's out = 1