Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF092617\
Data File : BF098827.D
Acq On : 26 Sep 2017 12:47
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Sep 27 00:59:02 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF092317.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Sep 23 13:50:58 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	105	0.00
2		1,4-Dioxane	0.600	0.581	3.2	100	-0.03
3		Pyridine	1.624	1.590	2.1	105	-0.02
4		n-Nitrosodimethylamine	0.688	0.669	2.8	102	-0.02
5	S	2-Fluorophenol	1.256	1.255	0.1	104	-0.01
6		Aniline	2.092	2.047	2.2	103	0.00
7	S	Phenol-d6	1.550	1.540	0.6	104	0.00
8		2-Chlorophenol	1.423	1.399	1.7	104	-0.01
9		Benzaldehyde	0.899	0.831	7.6	99	0.00
10	C	Phenol	1.733	1.677	3.2	103	0.00
11		bis(2-Chloroethyl)ether	1.348	1.319	2.2	104	0.00
12		1,3-Dichlorobenzene	1.490	1.466	1.6	105	0.00
13	C	1,4-Dichlorobenzene	1.516	1.472	2.9	104	0.00
14		1,2-Dichlorobenzene	1.401	1.366	2.5	103	0.00
15		Benzyl Alcohol	1.137	1.116	1.8	103	0.00
16		2,2'-oxybis(1-Chloropropane	2.099	2.018	3.9	104	0.00
17		2-Methylphenol	1.164	1.131	2.8	103	0.00
18		Hexachloroethane	0.545	0.527	3.3	103	-0.01
19	P	n-Nitroso-di-n-propylamine	0.990	0.928	6.3	103	0.00
20		3+4-Methylphenols	1.473	1.413	4.1	101	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	104	0.00
22		Acetophenone	0.485	0.458	5.6	102	0.00
23	S	Nitrobenzene-d5	0.312	0.311	0.3	103	0.00
24		Nitrobenzene	0.354	0.348	1.7	103	0.00
25		Isophorone	0.645	0.623	3.4	104	0.00
26	C	2-Nitrophenol	0.170	0.182	-7.1	108	-0.01
27		2,4-Dimethylphenol	0.321	0.309	3.7	103	0.00
28		bis(2-Chloroethoxy)methane	0.402	0.392	2.5	105	0.00
29	C	2,4-Dichlorophenol	0.276	0.274	0.7	105	0.00
30		1,2,4-Trichlorobenzene	0.276	0.272	1.4	104	0.00
31		Naphthalene	0.939	0.897	4.5	103	0.00
32		Benzoic acid	0.264	0.264	0.0	101	0.00
33		4-Chloroaniline	0.392	0.386	1.5	104	-0.01
34	C	Hexachlorobutadiene	0.142	0.141	0.7	107	0.00
35		Caprolactam	0.088	0.086	2.3	105	0.00
36	C	4-Chloro-3-methylphenol	0.310	0.301	2.9	103	0.00
37		2-Methylnaphthalene	0.611	0.587	3.9	103	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	104	0.00
39		1,2,4,5-Tetrachlorobenzene	0.596	0.585	1.8	105	0.00
40	P	Hexachlorocyclopentadiene	0.273	0.264	3.3	99	0.00
41	S	2,4,6-Tribromophenol	0.164	0.168	-2.4	104	0.00
42	C	2,4,6-Trichlorophenol	0.423	0.411	2.8	103	0.00
43		2,4,5-Trichlorophenol	0.422	0.423	-0.2	104	0.00
44	S	2-Fluorobiphenyl	1.198	1.177	1.8	102	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.719	1.659	3.5	102	0.00
46		2-Chloronaphthalene	1.291	1.258	2.6	104	-0.01
47		2-Nitroaniline	0.395	0.397	-0.5	103	0.00
48		Acenaphthylene	1.976	1.917	3.0	103	-0.01
49		Dimethylphthalate	1.509	1.465	2.9	104	0.00
50		2,6-Dinitrotoluene	0.329	0.333	-1.2	103	0.00
51	C	Acenaphthene	1.251	1.221	2.4	104	0.00
52		3-Nitroaniline	0.383	0.389	-1.6	102	0.00
53	P	2,4-Dinitrophenol	0.148	0.155	-4.7	104	0.00
54		Dibenzofuran	1.666	1.602	3.8	102	0.00
55	P	4-Nitrophenol	0.324	0.327	-0.9	103	0.00
56		2,4-Dinitrotoluene	0.426	0.439	-3.1	104	0.00
57		Fluorene	1.339	1.274	4.9	102	0.00
58		2,3,4,6-Tetrachlorophenol	0.291	0.290	0.3	104	0.00
59		Diethylphthalate	1.475	1.416	4.0	102	-0.01
60		4-Chlorophenyl-phenylether	0.602	0.578	4.0	103	0.00
61		4-Nitroaniline	0.370	0.375	-1.4	103	0.00
62		Azobenzene	1.356	1.293	4.6	102	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
64		4,6-Dinitro-2-methylphenol	0.122	0.129	-5.7	106	0.00
65	c	n-Nitrosodiphenylamine	0.693	0.689	0.6	104	0.00
66		4-Bromophenyl-phenylether	0.196	0.195	0.5	103	-0.01
67		Hexachlorobenzene	0.209	0.209	0.0	104	0.00
68		Atrazine	0.208	0.199	4.3	98	0.00
69	C	Pentachlorophenol	0.130	0.129	0.8	97	0.00
70		Phenanthrene	1.071	1.038	3.1	102	0.00
71		Anthracene	1.095	1.078	1.6	102	-0.01
72		Carbazole	1.073	1.042	2.9	100	0.00
73		Di-n-butylphthalate	1.291	1.261	2.3	100	-0.01
74	C	Fluoranthene	1.084	1.040	4.1	100	-0.01
75	I	Chrysene-d12	1.000	1.000	0.0	102	0.00
76		Benzidine	0.740	0.763	-3.1	106	-0.01
77		Pyrene	1.525	1.483	2.8	101	0.00
78	S	Terphenyl-d14	0.879	0.870	1.0	102	-0.01
79		Butylbenzylphthalate	0.717	0.710	1.0	100	-0.01
80		Benzo(a)anthracene	1.211	1.187	2.0	103	0.00
81		3,3'-Dichlorobenzidine	0.444	0.442	0.5	102	0.00
82		Chrysene	1.153	1.116	3.2	101	0.00
83		Bis(2-ethylhexyl)phthalate	0.987	0.981	0.6	101	0.00
84	c	Di-n-octyl phthalate	1.589	1.596	-0.4	105	0.00
85		Indeno(1,2,3-cd)pyrene	0.922	0.966	-4.8	118	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	111	0.00
87		Benzo(b)fluoranthene	1.230	1.233	-0.2	109	-0.01

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88	Benzo(k)fluoranthene	1.215	1.136	6.5	105	0.00
89 C	Benzo(a)pyrene	1.129	1.102	2.4	109	-0.01
90	Dibenzo(a,h)anthracene	0.903	0.914	-1.2	116	0.00
91	Benzo(g,h,i)perylene	0.893	0.920	-3.0	117	0.00

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