Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\Data\BM052417\
Data File : BM010172.D
Acq On : 24 May 2017 10:01
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: May 25 05:22:40 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM051917.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat May 20 04:25:49 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	63	-0.02
2		1,4-Dioxane	0.509	0.493	3.1	63	-0.02
3		Pyridine	1.362	1.374	-0.9	62	-0.03
4		n-Nitrosodimethylamine	0.493	0.607	-23.1	83	-0.02
5	S	2-Fluorophenol	1.108	1.052	5.1	61	-0.02
6		Aniline	1.766	1.691	4.2	60	-0.02
7	S	Phenol-d6	1.425	1.369	3.9	61	-0.03
8		2-Chlorophenol	1.205	1.132	6.1	61	-0.02
9		Benzaldehyde	0.887	0.877	1.1	61	-0.02
10	C	Phenol	1.502	1.448	3.6	62	-0.03
11		bis(2-Chloroethyl)ether	1.119	1.062	5.1	59	-0.02
12		1,3-Dichlorobenzene	1.531	1.483	3.1	62	-0.02
13	C	1,4-Dichlorobenzene	1.573	1.507	4.2	61	-0.02
14		1,2-Dichlorobenzene	1.510	1.459	3.4	62	-0.02
15		Benzyl Alcohol	1.078	1.105	-2.5	63	-0.02
16		2,2'-oxybis(1-Chloropropane	1.101	0.905	17.8	51	-0.04
17		2-Methylphenol	1.069	1.006	5.9	58	-0.03
18		Hexachloroethane	0.510	0.510	0.0	64	-0.02
19	P	n-Nitroso-di-n-propylamine	1.006	1.023	-1.7	63	-0.03
20		3+4-Methylphenols	1.443	1.384	4.1	60	-0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	57	-0.02
22		Acetophenone	0.513	0.535	-4.3	61	-0.03
23	S	Nitrobenzene-d5	0.371	0.432	-16.4	65	-0.03
24		Nitrobenzene	0.382	0.440	-15.2	65	-0.02
25		Isophorone	0.641	0.694	-8.3	61	-0.04
26	C	2-Nitrophenol	0.161	0.176	-9.3	61	-0.02
27		2,4-Dimethylphenol	0.291	0.294	-1.0	58	-0.02
28		bis(2-Chloroethoxy)methane	0.407	0.411	-1.0	57	-0.02
29	C	2,4-Dichlorophenol	0.335	0.370	-10.4	62	-0.03
30		1,2,4-Trichlorobenzene	0.433	0.457	-5.5	62	-0.02
31		Naphthalene	1.070	1.051	1.8	57	-0.03
32		Benzoic acid	0.198	0.204	-3.0	58	-0.04
33		4-Chloroaniline	0.412	0.417	-1.2	56	-0.02
34	C	Hexachlorobutadiene	0.289	0.331	-14.5	66	-0.02
35		Caprolactam	0.096	0.098	-2.1	55	-0.05
36	C	4-Chloro-3-methylphenol	0.362	0.382	-5.5	59	-0.02
37		2-Methylnaphthalene	0.770	0.791	-2.7	59	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	58	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.769	0.824	-7.2	64	-0.02
40	P	Hexachlorocyclopentadiene	0.444	0.480	-8.1	63	-0.02
41	S	2,4,6-Tribromophenol	0.304	0.326	-7.2	62	-0.02
42	C	2,4,6-Trichlorophenol	0.453	0.507	-11.9	64	-0.02
43		2,4,5-Trichlorophenol	0.501	0.557	-11.2	63	-0.02
44	S	2-Fluorobiphenyl	1.556	1.636	-5.1	62	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.645	1.690	-2.7	61	-0.02
46		2-Chloronaphthalene	1.332	1.363	-2.3	61	-0.02
47		2-Nitroaniline	0.342	0.392	-14.6	67	-0.02
48		Acenaphthylene	1.969	2.009	-2.0	59	-0.02
49		Dimethylphthalate	1.784	1.830	-2.6	60	-0.02
50		2,6-Dinitrotoluene	0.339	0.356	-5.0	59	-0.02
51	C	Acenaphthene	1.224	1.242	-1.5	59	-0.02
52		3-Nitroaniline	0.315	0.311	1.3	58	-0.02
53	P	2,4-Dinitrophenol	0.195	0.212	-8.7	65	-0.02
54		Dibenzofuran	1.925	1.966	-2.1	60	-0.02
55	P	4-Nitrophenol	0.253	0.245	3.2	56	-0.03
56		2,4-Dinitrotoluene	0.479	0.502	-4.8	61	-0.02
57		Fluorene	1.674	1.741	-4.0	62	-0.02
58		2,3,4,6-Tetrachlorophenol	0.426	0.477	-12.0	64	-0.02
59		Diethylphthalate	1.687	1.757	-4.1	61	-0.03
60		4-Chlorophenyl-phenylether	0.940	1.003	-6.7	64	-0.02
61		4-Nitroaniline	0.320	0.314	1.9	58	-0.02
62		Azobenzene	1.240	1.476	-19.0	69	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	62	-0.02
64		4,6-Dinitro-2-methylphenol	0.122	0.130	-6.6	64	-0.02
65	c	n-Nitrosodiphenylamine	0.599	0.613	-2.3	62	-0.02
66		4-Bromophenyl-phenylether	0.256	0.269	-5.1	65	-0.02
67		Hexachlorobenzene	0.311	0.306	1.6	61	-0.02
68		Atrazine	0.227	0.242	-6.6	63	-0.03
69	C	Pentachlorophenol	0.171	0.186	-8.8	64	-0.02
70		Phenanthrene	1.119	1.139	-1.8	63	-0.02
71		Anthracene	1.110	1.167	-5.1	64	-0.02
72		Carbazole	0.983	1.003	-2.0	62	-0.02
73		Di-n-butylphthalate	1.166	1.183	-1.5	61	-0.02
74	C	Fluoranthene	1.453	1.520	-4.6	65	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	67	-0.02
76		Benzidine	0.369	0.536	-45.3#	88	-0.02
77		Pyrene	1.072	1.051	2.0	65	-0.02
78	S	Terphenyl-d14	0.880	0.914	-3.9	68	-0.02
79		Butylbenzylphthalate	0.397	0.384	3.3	63	-0.02
80		Benzo(a)anthracene	1.102	1.121	-1.7	68	-0.02
81		3,3'-Dichlorobenzidine	0.443	0.459	-3.6	68	-0.02
82		Chrysene	1.045	1.063	-1.7	68	-0.02
83		Bis(2-ethylhexyl)phthalate	0.594	0.580	2.4	65	-0.03
84	c	Di-n-octyl phthalate	1.063	1.065	-0.2	66	-0.04
85		Indeno(1,2,3-cd)pyrene	1.533	1.659	-8.2	74	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	68	-0.03
87		Benzo(b)fluoranthene	1.170	1.214	-3.8	69	-0.03

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88	Benzo(k)fluoranthene	1.147	1.155	-0.7	69	-0.03
89 C	Benzo(a)pyrene	1.115	1.144	-2.6	70	-0.04
90	Dibenzo(a,h)anthracene	1.247	1.324	-6.2	73	-0.05
91	Benzo(g,h,i)perylene	1.222	1.290	-5.6	72	-0.06

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