Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF100317\
Data File : BF099036.D
Acq On : 3 Oct 2017 17:53
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Oct 04 02:19:27 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	76	-0.05
2		1,4-Dioxane	0.600	0.605	-0.8	76	-0.11
3		Pyridine	1.624	1.586	2.3	76	-0.10
4		n-Nitrosodimethylamine	0.688	0.641	6.8	71	-0.11
5	S	2-Fluorophenol	1.256	1.316	-4.8	80	-0.05
6		Aniline	2.092	2.139	-2.2	79	-0.05
7	S	Phenol-d6	1.550	1.625	-4.8	80	-0.04
8		2-Chlorophenol	1.423	1.485	-4.4	81	-0.05
9		Benzaldehyde	0.899	0.843	6.2	73	-0.05
10	C	Phenol	1.733	1.797	-3.7	80	-0.04
11		bis(2-Chloroethyl)ether	1.348	1.385	-2.7	79	-0.05
12		1,3-Dichlorobenzene	1.490	1.526	-2.4	79	-0.05
13	C	1,4-Dichlorobenzene	1.516	1.577	-4.0	81	-0.05
14		1,2-Dichlorobenzene	1.401	1.454	-3.8	80	-0.05
15		Benzyl Alcohol	1.137	1.136	0.1	76	-0.05
16		2,2'-oxybis(1-Chloropropane	2.099	2.043	2.7	76	-0.05
17		2-Methylphenol	1.164	1.194	-2.6	79	-0.04
18		Hexachloroethane	0.545	0.552	-1.3	79	-0.05
19	P	n-Nitroso-di-n-propylamine	0.990	0.951	3.9	77	-0.05
20		3+4-Methylphenols	1.473	1.474	-0.1	77	-0.05
21	I	Naphthalene-d8	1.000	1.000	0.0	74	-0.05
22		Acetophenone	0.485	0.483	0.4	77	-0.05
23	S	Nitrobenzene-d5	0.312	0.335	-7.4	79	-0.05
24		Nitrobenzene	0.354	0.367	-3.7	77	-0.04
25		Isophorone	0.645	0.651	-0.9	78	-0.05
26	C	2-Nitrophenol	0.170	0.200	-17.6	84	-0.05
27		2,4-Dimethylphenol	0.321	0.331	-3.1	78	-0.05
28		bis(2-Chloroethoxy)methane	0.402	0.414	-3.0	79	-0.05
29	C	2,4-Dichlorophenol	0.276	0.294	-6.5	80	-0.05
30		1,2,4-Trichlorobenzene	0.276	0.294	-6.5	80	-0.05
31		Naphthalene	0.939	0.979	-4.3	80	-0.05
32		Benzoic acid	0.264	0.262	0.8	72	-0.04
33		4-Chloroaniline	0.392	0.411	-4.8	79	-0.05
34	C	Hexachlorobutadiene	0.142	0.148	-4.2	80	-0.05
35		Caprolactam	0.088	0.088	0.0	77	-0.05
36	C	4-Chloro-3-methylphenol	0.310	0.317	-2.3	77	-0.04
37		2-Methylnaphthalene	0.611	0.631	-3.3	79	-0.05
38	I	Acenaphthene-d10	1.000	1.000	0.0	74	-0.05
39		1,2,4,5-Tetrachlorobenzene	0.596	0.641	-7.6	81	-0.05
40	P	Hexachlorocyclopentadiene	0.273	0.272	0.4	71	-0.05
41	S	2,4,6-Tribromophenol	0.164	0.174	-6.1	76	-0.05
42	C	2,4,6-Trichlorophenol	0.423	0.439	-3.8	77	-0.05
43		2,4,5-Trichlorophenol	0.422	0.444	-5.2	77	-0.04
44	S	2-Fluorobiphenyl	1.198	1.303	-8.8	80	-0.05

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.719	1.812	-5.4	79	-0.05
46		2-Chloronaphthalene	1.291	1.364	-5.7	79	-0.05
47		2-Nitroaniline	0.395	0.410	-3.8	75	-0.04
48		Acenaphthylene	1.976	2.070	-4.8	79	-0.05
49		Dimethylphthalate	1.509	1.544	-2.3	77	-0.05
50		2,6-Dinitrotoluene	0.329	0.349	-6.1	76	-0.04
51	C	Acenaphthene	1.251	1.320	-5.5	79	-0.05
52		3-Nitroaniline	0.383	0.413	-7.8	76	-0.04
53	P	2,4-Dinitrophenol	0.148	0.172	-16.2	82	-0.03
54		Dibenzofuran	1.666	1.743	-4.6	78	-0.05
55	P	4-Nitrophenol	0.324	0.321	0.9	71	-0.04
56		2,4-Dinitrotoluene	0.426	0.462	-8.5	77	-0.04
57		Fluorene	1.339	1.396	-4.3	79	-0.05
58		2,3,4,6-Tetrachlorophenol	0.291	0.301	-3.4	76	-0.05
59		Diethylphthalate	1.475	1.493	-1.2	76	-0.05
60		4-Chlorophenyl-phenylether	0.602	0.626	-4.0	78	-0.05
61		4-Nitroaniline	0.370	0.404	-9.2	78	-0.05
62		Azobenzene	1.356	1.340	1.2	74	-0.05
63	I	Phenanthrene-d10	1.000	1.000	0.0	70	-0.05
64		4,6-Dinitro-2-methylphenol	0.122	0.142	-16.4	81	-0.04
65	c	n-Nitrosodiphenylamine	0.693	0.734	-5.9	77	-0.05
66		4-Bromophenyl-phenylether	0.196	0.213	-8.7	78	-0.05
67		Hexachlorobenzene	0.209	0.221	-5.7	77	-0.05
68		Atrazine	0.208	0.220	-5.8	76	-0.04
69	C	Pentachlorophenol	0.130	0.124	4.6	65	-0.05
70		Phenanthrene	1.071	1.111	-3.7	76	-0.05
71		Anthracene	1.095	1.126	-2.8	75	-0.05
72		Carbazole	1.073	1.093	-1.9	73	-0.05
73		Di-n-butylphthalate	1.291	1.320	-2.2	73	-0.05
74	C	Fluoranthene	1.084	1.071	1.2	72	-0.05
75	I	Chrysene-d12	1.000	1.000	0.0	59	-0.05
76		Benzidine	0.740	0.875	-18.2	71	-0.05
77		Pyrene	1.525	1.802	-18.2	71	-0.05
78	S	Terphenyl-d14	0.879	1.060	-20.6	72	-0.05
79		Butylbenzylphthalate	0.717	0.788	-9.9	64	-0.05
80		Benzo(a)anthracene	1.211	1.275	-5.3	64	-0.05
81		3,3'-Dichlorobenzidine	0.444	0.471	-6.1	63	-0.05
82		Chrysene	1.153	1.212	-5.1	64	-0.05
83		Bis(2-ethylhexyl)phthalate	0.987	1.015	-2.8	61	-0.05
84	c	Di-n-octyl phthalate	1.589	1.526	4.0	58	-0.05
85		Indeno(1,2,3-cd)pyrene	0.922	1.043	-13.1	74	-0.09
86	I	Perylene-d12	1.000	1.000	0.0	60	-0.06
87		Benzo(b)fluoranthene	1.230	1.276	-3.7	62	-0.06

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.215	1.245	-2.5	63	-0.06
89 C	Benzo(a)pyrene	1.129	1.194	-5.8	64	-0.06
90	Dibenzo(a,h)anthracene	0.903	1.062	-17.6	73	-0.09
91	Benzo(g,h,i)perylene	0.893	1.084	-21.4	75	-0.09

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