Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF071116\
Data File : BF088898.D
Acq On : 11 Jul 2016 14:25
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 11 18:15:12 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jul 11 18:13:32 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	56	0.00
2		1,4-Dioxane	0.662	0.613	7.4	54	0.00
3		Pyridine	1.920	1.681	12.4	51	0.00
4		n-Nitrosodimethylamine	0.733	0.615	16.1	49#	0.00
5	S	2-Fluorophenol	1.334	1.281	4.0	54	0.00
6		Aniline	2.513	2.341	6.8	52	0.00
7	S	Phenol-d6	1.724	1.609	6.7	55	0.00
8		2-Chlorophenol	1.434	1.399	2.4	59	0.00
9		Benzaldehyde	1.168	0.955	18.2	49#	0.00
10	C	Phenol	1.968	1.758	10.7	50	0.00
11		bis(2-Chloroethyl)ether	1.468	1.469	-0.1	60	0.00
12		1,3-Dichlorobenzene	1.578	1.621	-2.7	63	0.00
13	C	1,4-Dichlorobenzene	1.567	1.700	-8.5	65	0.00
14		1,2-Dichlorobenzene	1.466	1.417	3.3	61	0.00
15		Benzyl Alcohol	1.269	1.154	9.1	50#	0.00
16		2,2'-oxybis(1-Chloropropane	2.125	1.720	19.1	46#	0.00
17		2-Methylphenol	1.170	1.205	-3.0	58	0.00
18		Hexachloroethane	0.606	0.604	0.3	58	0.00
19	P	n-Nitroso-di-n-propylamine	1.158	1.132	2.2	64	0.00
20		3+4-Methylphenols	1.404	1.320	6.0	57	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	58	0.00
22		Acetophenone	0.485	0.499	-2.9	61	0.00
23	S	Nitrobenzene-d5	0.382	0.333	12.8	53	0.00
24		Nitrobenzene	0.385	0.396	-2.9	62	0.00
25		Isophorone	0.707	0.629	11.0	54	0.00
26	C	2-Nitrophenol	0.158	0.163	-3.2	64	0.00
27		2,4-Dimethylphenol	0.329	0.310	5.8	54	0.00
28		bis(2-Chloroethoxy)methane	0.460	0.446	3.0	61	0.00
29	C	2,4-Dichlorophenol	0.266	0.276	-3.8	64	0.00
30		1,2,4-Trichlorobenzene	0.291	0.299	-2.7	60	0.00
31		Naphthalene	0.986	1.065	-8.0	62	0.00
32		Benzoic acid	0.239	0.242	-1.3	59	0.00
33		4-Chloroaniline	0.439	0.396	9.8	53	0.00
34	C	Hexachlorobutadiene	0.156	0.173	-10.9	64	0.00
35		Caprolactam	0.090	0.086	4.4	54	0.00
36	C	4-Chloro-3-methylphenol	0.314	0.317	-1.0	59	0.00
37		2-Methylnaphthalene	0.635	0.589	7.2	61	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	60	0.00
39		1,2,4,5-Tetrachlorobenzene	0.665	0.756	-13.7	66	0.00
40	P	Hexachlorocyclopentadiene	0.327	0.289	11.6	53	0.00
41	S	2,4,6-Tribromophenol	0.186	0.231	-24.2	71	0.00
42	C	2,4,6-Trichlorophenol	0.439	0.440	-0.2	65	0.00
43		2,4,5-Trichlorophenol	0.377	0.408	-8.2	63	0.00
44	S	2-Fluorobiphenyl	1.266	1.276	-0.8	67	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.656	1.801	-8.8	64	0.00
46		2-Chloronaphthalene	1.300	1.394	-7.2	62	0.00
47		2-Nitroaniline	0.461	0.399	13.4	47#	0.00
48		Acenaphthylene	2.069	2.237	-8.1	64	0.00
49		Dimethylphthalate	1.425	1.560	-9.5	71	0.00
50		2,6-Dinitrotoluene	0.316	0.365	-15.5	74	0.00
51	C	Acenaphthene	1.268	1.377	-8.6	68	0.00
52		3-Nitroaniline	0.401	0.352	12.2	46#	0.00
53	P	2,4-Dinitrophenol	0.139	0.146	-5.0	64	0.00
54		Dibenzofuran	1.660	1.770	-6.6	64	0.00
55	P	4-Nitrophenol	0.305	0.299	2.0	53	0.00
56		2,4-Dinitrotoluene	0.413	0.468	-13.3	71	0.00
57		Fluorene	1.279	1.448	-13.2	65	0.00
58		2,3,4,6-Tetrachlorophenol	0.292	0.298	-2.1	59	0.00
59		Diethylphthalate	1.430	1.499	-4.8	63	0.00
60		4-Chlorophenyl-phenylether	0.594	0.573	3.5	63	0.00
61		4-Nitroaniline	0.386	0.410	-6.2	69	0.00
62		Azobenzene	1.631	1.528	6.3	59	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	63	0.00
64		4,6-Dinitro-2-methylphenol	0.107	0.111	-3.7	60	0.00
65	c	n-Nitrosodiphenylamine	0.677	0.694	-2.5	62	0.00
66		4-Bromophenyl-phenylether	0.202	0.212	-5.0	68	0.00
67		Hexachlorobenzene	0.223	0.215	3.6	61	0.00
68		Atrazine	0.211	0.183	13.3	53	0.00
69	C	Pentachlorophenol	0.132	0.122	7.6	55	0.00
70		Phenanthrene	1.093	0.992	9.2	61	0.00
71		Anthracene	1.087	1.143	-5.2	66	0.00
72		Carbazole	1.023	0.892	12.8	61	0.00
73		Di-n-butylphthalate	1.190	1.155	2.9	65	0.00
74	C	Fluoranthene	1.108	1.043	5.9	57	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	68	0.00
76		Benzidine	1.011	1.098	-8.6	72	0.00
77		Pyrene	1.696	1.446	14.7	55	0.00
78	S	Terphenyl-d14	0.898	0.901	-0.3	70	0.00
79		Butylbenzylphthalate	0.756	0.660	12.7	60	0.00
80		Benzo(a)anthracene	1.288	1.256	2.5	66	0.00
81		3,3'-Dichlorobenzidine	0.484	0.501	-3.5	73	0.00
82		Chrysene	1.274	1.262	0.9	68	0.00
83		Bis(2-ethylhexyl)phthalate	0.864	0.877	-1.5	65	0.00
84	c	Di-n-octyl phthalate	1.467	1.599	-9.0	71	0.00
85		Indeno(1,2,3-cd)pyrene	0.980	0.913	6.8	66	0.00
86	I	Perylene-d12	1.000	1.000	0.0	65	0.00
87		Benzo(b)fluoranthene	1.255	1.162	7.4	63	0.00

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88	Benzo(k)fluoranthene	1.092	1.266	-15.9	75	0.00
89 C	Benzo(a)pyrene	1.062	1.115	-5.0	68	0.00
90	Dibenzo(a,h)anthracene	0.887	0.900	-1.5	67	0.00
91	Benzo(g,h,i)perylene	0.918	0.896	2.4	64	0.00

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