Evaluate Continuing Calibration Report

Data Path : V:\HPCHEM1\BNA F\DATA\BF070816\
Data File : BF088856.D
Acq On : 9 Jul 2016 6:36
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jul 09 07:04:02 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF063016.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 08 18:51:20 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	75	0.00
2		1,4-Dioxane	0.662	0.580	12.4	67	0.00
3		Pyridine	1.920	1.637	14.7	66	0.00
4		n-Nitrosodimethylamine	0.733	0.594	19.0	63	0.00
5	S	2-Fluorophenol	1.334	1.170	12.3	65	-0.01
6		Aniline	2.513	2.275	9.5	68	0.00
7	S	Phenol-d6	1.724	1.547	10.3	70	0.00
8		2-Chlorophenol	1.434	1.324	7.7	74	0.00
9		Benzaldehyde	1.168	0.933	20.1	64	0.00
10	C	Phenol	1.968	1.631	17.1	62	0.00
11		bis(2-Chloroethyl)ether	1.468	1.393	5.1	76	0.00
12		1,3-Dichlorobenzene	1.578	1.618	-2.5	84	0.00
13	C	1,4-Dichlorobenzene	1.567	1.624	-3.6	83	0.00
14		1,2-Dichlorobenzene	1.466	1.344	8.3	77	0.00
15		Benzyl Alcohol	1.269	1.102	13.2	63	0.00
16		2,2'-oxybis(1-Chloropropane	2.125	1.638	22.9	59	0.00
17		2-Methylphenol	1.170	1.147	2.0	74	0.01
18		Hexachloroethane	0.606	0.602	0.7	76	0.00
19	P	n-Nitroso-di-n-propylamine	1.158	1.107	4.4	83	0.00
20		3+4-Methylphenols	1.404	1.233	12.2	71	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	73	0.00
22		Acetophenone	0.485	0.498	-2.7	77	0.00
23	S	Nitrobenzene-d5	0.382	0.335	12.3	68	0.00
24		Nitrobenzene	0.385	0.400	-3.9	80	0.00
25		Isophorone	0.707	0.651	7.9	70	0.00
26	C	2-Nitrophenol	0.158	0.157	0.6	79	0.00
27		2,4-Dimethylphenol	0.329	0.290	11.9	64	0.00
28		bis(2-Chloroethoxy)methane	0.460	0.458	0.4	80	0.00
29	C	2,4-Dichlorophenol	0.266	0.282	-6.0	82	0.00
30		1,2,4-Trichlorobenzene	0.291	0.315	-8.2	79	0.00
31		Naphthalene	0.986	1.075	-9.0	79	0.00
32		Benzoic acid	0.239	0.197	17.6	61	0.00
33		4-Chloroaniline	0.439	0.414	5.7	70	0.00
34	C	Hexachlorobutadiene	0.156	0.161	-3.2	76	0.00
35		Caprolactam	0.090	0.090	0.0	71	0.01
36	C	4-Chloro-3-methylphenol	0.314	0.340	-8.3	80	0.00
37		2-Methylnaphthalene	0.635	0.645	-1.6	84	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	74	0.00
39		1,2,4,5-Tetrachlorobenzene	0.665	0.768	-15.5	83	0.00
40	P	Hexachlorocyclopentadiene	0.327	0.274	16.2	63	0.00
41	S	2,4,6-Tribromophenol	0.186	0.217	-16.7	83	0.00
42	C	2,4,6-Trichlorophenol	0.439	0.406	7.5	75	0.00
43		2,4,5-Trichlorophenol	0.377	0.436	-15.6	83	0.00
44	S	2-Fluorobiphenyl	1.266	1.249	1.3	81	0.00

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ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

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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)
45		1,1'-Biphenyl	1.656	1.875	-13.2	82	0.00
46		2-Chloronaphthalene	1.300	1.394	-7.2	78	0.00
47		2-Nitroaniline	0.461	0.455	1.3	66	0.00
48		Acenaphthylene	2.069	2.232	-7.9	80	0.00
49		Dimethylphthalate	1.425	1.481	-3.9	84	0.00
50		2,6-Dinitrotoluene	0.316	0.326	-3.2	82	-0.01
51	C	Acenaphthene	1.268	1.395	-10.0	86	0.00
52		3-Nitroaniline	0.401	0.380	5.2	62	0.00
53	P	2,4-Dinitrophenol	0.139	0.104	25.2#	57	0.00
54		Dibenzofuran	1.660	1.716	-3.4	78	0.00
55	P	4-Nitrophenol	0.305	0.246	19.3	55	0.00
56		2,4-Dinitrotoluene	0.413	0.431	-4.4	82	0.00
57		Fluorene	1.279	1.276	0.2	71	-0.01
58		2,3,4,6-Tetrachlorophenol	0.292	0.311	-6.5	77	0.00
59		Diethylphthalate	1.430	1.415	1.0	74	0.00
60		4-Chlorophenyl-phenylether	0.594	0.608	-2.4	83	0.00
61		4-Nitroaniline	0.386	0.410	-6.2	86	0.00
62		Azobenzene	1.631	1.639	-0.5	78	0.00
63	I	Phenanthrene-d10	1.000	1.000	0.0	79	0.00
64		4,6-Dinitro-2-methylphenol	0.107	0.087	18.7	58	0.00
65	c	n-Nitrosodiphenylamine	0.677	0.669	1.2	75	0.00
66		4-Bromophenyl-phenylether	0.202	0.211	-4.5	85	0.00
67		Hexachlorobenzene	0.223	0.204	8.5	72	0.00
68		Atrazine	0.211	0.210	0.5	75	0.00
69	C	Pentachlorophenol	0.132	0.116	12.1	65	0.00
70		Phenanthrene	1.093	1.036	5.2	79	0.00
71		Anthracene	1.087	1.038	4.5	75	0.00
72		Carbazole	1.023	0.900	12.0	77	0.00
73		Di-n-butylphthalate	1.190	1.249	-5.0	87	0.00
74	C	Fluoranthene	1.108	0.955	13.8	65	0.00
75	I	Chrysene-d12	1.000	1.000	0.0	60	-0.01
76		Benzidine	1.011	0.930	8.0	53	0.00
77		Pyrene	1.696	1.985	-17.0	66	0.00
78	S	Terphenyl-d14	0.898	1.033	-15.0	70	0.00
79		Butylbenzylphthalate	0.756	0.909	-20.2	72	0.00
80		Benzo(a)anthracene	1.288	1.343	-4.3	62	0.00
81		3,3'-Dichlorobenzidine	0.484	0.467	3.5	60	0.00
82		Chrysene	1.274	1.394	-9.4	66	0.00
83		Bis(2-ethylhexyl)phthalate	0.864	1.070	-23.8	70	0.00
84	c	Di-n-octyl phthalate	1.467	1.671	-13.9	65	0.00
85		Indeno(1,2,3-cd)pyrene	0.980	1.475	-50.5#	94	0.00
86	I	Perylene-d12	1.000	1.000	0.0	72	0.00
87		Benzo(b)fluoranthene	1.255	1.049	16.4	63	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.092	1.137	-4.1	75	0.00
89 C	Benzo(a)pyrene	1.062	1.038	2.3	70	0.00
90	Dibenzo(a,h)anthracene	0.887	1.132	-27.6#	94	0.00
91	Benzo(g,h,i)perylene	0.918	1.180	-28.5#	94	0.01

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