Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061116\
Data File : BF087921.D
Acq On : 12 Jun 2016 00:28
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 12 05:56:04 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 07 18:51:55 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	81	0.01
2		1,4-Dioxane	0.568	0.600	-5.6	88	0.02
3		Pyridine	1.342	1.820	-35.6#	108	0.01
4		n-Nitrosodimethylamine	0.568	0.659	-16.0	94	0.02
5	S	2-Fluorophenol	1.170	1.203	-2.8	85	0.00
6		Aniline	2.018	2.341	-16.0	95	0.01
7	S	Phenol-d6	1.418	1.693	-19.4	101	0.00
8		2-Chlorophenol	1.385	1.331	3.9	80	0.00
9		Benzaldehyde	0.923	0.922	0.1	87	0.01
10	C	Phenol	1.665	1.980	-18.9	115	0.01
11		bis(2-Chloroethyl)ether	1.243	1.286	-3.5	90	0.01
12		1,3-Dichlorobenzene	1.486	1.432	3.6	79	0.01
13	C	1,4-Dichlorobenzene	1.497	1.493	0.3	85	0.01
14		1,2-Dichlorobenzene	1.361	1.246	8.4	73	0.00
15		Benzyl Alcohol	1.109	0.992	10.6	70	0.01
16		2,2'-oxybis(1-Chloropropane	1.680	1.669	0.7	80	0.01
17		2-Methylphenol	1.123	1.139	-1.4	80	0.00
18		Hexachloroethane	0.531	0.514	3.2	78	0.00
19	P	n-Nitroso-di-n-propylamine	0.907	0.977	-7.7	95	0.01
20		3+4-Methylphenols	1.310	1.216	7.2	81	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	84	0.01
22		Acetophenone	0.447	0.425	4.9	82	0.01
23	S	Nitrobenzene-d5	0.343	0.327	4.7	78	0.01
24		Nitrobenzene	0.332	0.375	-13.0	103	0.01
25		Isophorone	0.634	0.632	0.3	82	0.01
26	C	2-Nitrophenol	0.178	0.190	-6.7	78	0.01
27		2,4-Dimethylphenol	0.327	0.313	4.3	78	0.00
28		bis(2-Chloroethoxy)methane	0.393	0.420	-6.9	87	0.01
29	C	2,4-Dichlorophenol	0.273	0.289	-5.9	87	0.01
30		1,2,4-Trichlorobenzene	0.297	0.281	5.4	82	0.01
31		Naphthalene	0.990	0.928	6.3	83	0.01
32		Benzoic acid	0.180	0.197	-9.4	78	0.01
33		4-Chloroaniline	0.442	0.464	-5.0	82	0.01
34	C	Hexachlorobutadiene	0.157	0.157	0.0	81	0.01
35		Caprolactam	0.090	0.088	2.2	74	0.01
36	C	4-Chloro-3-methylphenol	0.292	0.291	0.3	85	0.01
37		2-Methylnaphthalene	0.652	0.707	-8.4	86	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	97	0.01
39		1,2,4,5-Tetrachlorobenzene	0.583	0.452	22.5	87	0.01
40	P	Hexachlorocyclopentadiene	0.323	0.246	23.8	73	0.01
41	S	2,4,6-Tribromophenol	0.211	0.183	13.3	82	0.01
42	C	2,4,6-Trichlorophenol	0.400	0.381	4.8	90	0.01
43		2,4,5-Trichlorophenol	0.416	0.381	8.4	92	0.01
44	S	2-Fluorobiphenyl	1.502	1.179	21.5	92	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.613	1.427	11.5	93	0.00
46		2-Chloronaphthalene	1.294	1.168	9.7	95	0.00
47		2-Nitroaniline	0.382	0.399	-4.5	93	0.01
48		Acenaphthylene	2.059	1.962	4.7	94	0.01
49		Dimethylphthalate	1.449	1.256	13.3	93	0.01
50		2,6-Dinitrotoluene	0.332	0.290	12.7	94	0.01
51	C	Acenaphthene	1.284	1.237	3.7	94	0.01
52		3-Nitroaniline	0.383	0.377	1.6	92	0.01
53	P	2,4-Dinitrophenol	0.122	0.172	-41.0#	100	0.01
54		Dibenzofuran	1.769	1.778	-0.5	96	0.01
55	P	4-Nitrophenol	0.297	0.303	-2.0	87	0.01
56		2,4-Dinitrotoluene	0.426	0.391	8.2	96	0.00
57		Fluorene	1.378	1.221	11.4	96	0.01
58		2,3,4,6-Tetrachlorophenol	0.327	0.318	2.8	89	0.01
59		Diethylphthalate	1.466	1.290	12.0	90	0.01
60		4-Chlorophenyl-phenylether	0.588	0.495	15.8	89	0.01
61		4-Nitroaniline	0.363	0.379	-4.4	93	0.00
62		Azobenzene	1.450	1.369	5.6	91	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	94	0.01
64		4,6-Dinitro-2-methylphenol	0.108	0.129	-19.4	86	0.01
65	c	n-Nitrosodiphenylamine	0.664	0.678	-2.1	95	0.01
66		4-Bromophenyl-phenylether	0.215	0.213	0.9	91	0.01
67		Hexachlorobenzene	0.223	0.205	8.1	94	0.01
68		Atrazine	0.201	0.206	-2.5	88	0.01
69	C	Pentachlorophenol	0.118	0.125	-5.9	89	0.01
70		Phenanthrene	1.049	0.994	5.2	100	0.00
71		Anthracene	1.059	1.026	3.1	87	0.01
72		Carbazole	0.991	1.002	-1.1	105	0.00
73		Di-n-butylphthalate	1.254	1.109	11.6	83	0.01
74	C	Fluoranthene	1.108	1.012	8.7	84	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	97	0.01
76		Benzidine	0.554	0.662	-19.5	116	0.01
77		Pyrene	1.484	1.330	10.4	88	0.01
78	S	Terphenyl-d14	0.879	0.693	21.2	80	0.01
79		Butylbenzylphthalate	0.656	0.676	-3.0	96	0.01
80		Benzo(a)anthracene	1.140	1.001	12.2	93	0.01
81		3,3'-Dichlorobenzidine	0.306	0.338	-10.5	100	0.01
82		Chrysene	1.072	0.967	9.8	94	0.01
83		Bis(2-ethylhexyl)phthalate	0.799	0.721	9.8	90	0.01
84	c	Di-n-octyl phthalate	1.298	1.317	-1.5	100	0.01
85		Indeno(1,2,3-cd)pyrene	0.996	0.916	8.0	90	0.02
86	I	Perylene-d12	1.000	1.000	0.0	99	0.01
87		Benzo(b)fluoranthene	1.394	1.139	18.3	76	0.01

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88	Benzo(k)fluoranthene	1.052	1.206	-14.6	139	0.02
89 C	Benzo(a)pyrene	1.128	1.113	1.3	96	0.01
90	Dibenzo(a,h)anthracene	1.047	0.943	9.9	89	0.03
91	Benzo(g,h,i)perylene	1.078	0.977	9.4	88	0.02

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