Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\DATA\BF061216\
Data File : BF087949.D
Acq On : 12 Jun 2016 15:37
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 13 05:08:44 2016
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF060716.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 13 02:17:28 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	95	0.01
2		1,4-Dioxane	0.568	0.586	-3.2	101	0.03
3		Pyridine	1.342	1.588	-18.3	110	0.02
4		n-Nitrosodimethylamine	0.568	0.594	-4.6	100	0.03
5	S	2-Fluorophenol	1.170	1.169	0.1	96	0.00
6		Aniline	2.018	2.049	-1.5	97	0.01
7	S	Phenol-d6	1.418	1.493	-5.3	104	0.01
8		2-Chlorophenol	1.385	1.326	4.3	93	0.00
9		Benzaldehyde	0.923	0.823	10.8	91	0.01
10	C	Phenol	1.665	1.788	-7.4	121	0.01
11		bis(2-Chloroethyl)ether	1.243	1.279	-2.9	106	0.01
12		1,3-Dichlorobenzene	1.486	1.481	0.3	95	0.01
13	C	1,4-Dichlorobenzene	1.497	1.504	-0.5	101	0.01
14		1,2-Dichlorobenzene	1.361	1.266	7.0	87	0.00
15		Benzyl Alcohol	1.109	0.974	12.2	81	0.01
16		2,2'-oxybis(1-Chloropropane	1.680	1.636	2.6	92	0.00
17		2-Methylphenol	1.123	1.138	-1.3	94	0.00
18		Hexachloroethane	0.531	0.507	4.5	91	0.00
19	P	n-Nitroso-di-n-propylamine	0.907	0.953	-5.1	109	0.01
20		3+4-Methylphenols	1.310	1.128	13.9	88	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	100	0.01
22		Acetophenone	0.447	0.419	6.3	96	0.01
23	S	Nitrobenzene-d5	0.343	0.318	7.3	91	0.01
24		Nitrobenzene	0.332	0.348	-4.8	113	0.01
25		Isophorone	0.634	0.604	4.7	93	0.01
26	C	2-Nitrophenol	0.178	0.184	-3.4	90	0.01
27		2,4-Dimethylphenol	0.327	0.311	4.9	92	0.00
28		bis(2-Chloroethoxy)methane	0.393	0.411	-4.6	101	0.01
29	C	2,4-Dichlorophenol	0.273	0.287	-5.1	103	0.01
30		1,2,4-Trichlorobenzene	0.297	0.268	9.8	93	0.01
31		Naphthalene	0.990	0.887	10.4	94	0.01
32		Benzoic acid	0.180	0.177	1.7	84	0.01
33		4-Chloroaniline	0.442	0.443	-0.2	93	0.01
34	C	Hexachlorobutadiene	0.157	0.147	6.4	90	0.01
35		Caprolactam	0.090	0.093	-3.3	93	0.01
36	C	4-Chloro-3-methylphenol	0.292	0.276	5.5	96	0.01
37		2-Methylnaphthalene	0.652	0.636	2.5	92	0.01
38	I	Acenaphthene-d10	1.000	1.000	0.0	93	0.01
39		1,2,4,5-Tetrachlorobenzene	0.583	0.530	9.1	97	0.01
40	P	Hexachlorocyclopentadiene	0.323	0.269	16.7	76	0.01
41	S	2,4,6-Tribromophenol	0.211	0.183	13.3	78	0.01
42	C	2,4,6-Trichlorophenol	0.400	0.399	0.3	90	0.01
43		2,4,5-Trichlorophenol	0.416	0.399	4.1	92	0.00
44	S	2-Fluorobiphenyl	1.502	1.283	14.6	95	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.613	1.462	9.4	91	0.00
46		2-Chloronaphthalene	1.294	1.173	9.4	90	0.00
47		2-Nitroaniline	0.382	0.415	-8.6	93	0.01
48		Acenaphthylene	2.059	1.964	4.6	90	0.01
49		Dimethylphthalate	1.449	1.311	9.5	93	0.01
50		2,6-Dinitrotoluene	0.332	0.304	8.4	94	0.01
51	C	Acenaphthene	1.284	1.283	0.1	93	0.01
52		3-Nitroaniline	0.383	0.408	-6.5	95	0.01
53	P	2,4-Dinitrophenol	0.122	0.134	-9.8	74	0.01
54		Dibenzofuran	1.769	1.704	3.7	88	0.01
55	P	4-Nitrophenol	0.297	0.291	2.0	80	0.01
56		2,4-Dinitrotoluene	0.426	0.384	9.9	90	0.00
57		Fluorene	1.378	1.146	16.8	86	0.00
58		2,3,4,6-Tetrachlorophenol	0.327	0.339	-3.7	91	0.01
59		Diethylphthalate	1.466	1.395	4.8	92	0.01
60		4-Chlorophenyl-phenylether	0.588	0.539	8.3	93	0.01
61		4-Nitroaniline	0.363	0.372	-2.5	87	0.00
62		Azobenzene	1.450	1.505	-3.8	95	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	94	0.01
64		4,6-Dinitro-2-methylphenol	0.108	0.118	-9.3	79	0.01
65	c	n-Nitrosodiphenylamine	0.664	0.620	6.6	87	0.00
66		4-Bromophenyl-phenylether	0.215	0.210	2.3	89	0.01
67		Hexachlorobenzene	0.223	0.197	11.7	90	0.01
68		Atrazine	0.201	0.210	-4.5	90	0.01
69	C	Pentachlorophenol	0.118	0.117	0.8	84	0.01
70		Phenanthrene	1.049	0.906	13.6	91	0.00
71		Anthracene	1.059	1.080	-2.0	92	0.01
72		Carbazole	0.991	0.911	8.1	95	0.00
73		Di-n-butylphthalate	1.254	1.170	6.7	87	0.01
74	C	Fluoranthene	1.108	1.023	7.7	85	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	80	0.01
76		Benzidine	0.554	0.646	-16.6	93	0.01
77		Pyrene	1.484	1.642	-10.6	89	0.01
78	S	Terphenyl-d14	0.879	0.827	5.9	78	0.01
79		Butylbenzylphthalate	0.656	0.762	-16.2	89	0.01
80		Benzo(a)anthracene	1.140	1.133	0.6	86	0.00
81		3,3'-Dichlorobenzidine	0.306	0.365	-19.3	89	0.01
82		Chrysene	1.072	1.199	-11.8	95	0.01
83		Bis(2-ethylhexyl)phthalate	0.799	0.884	-10.6	91	0.01
84	c	Di-n-octyl phthalate	1.298	1.512	-16.5	94	0.01
85		Indeno(1,2,3-cd)pyrene	0.996	1.052	-5.6	85	0.02
86	I	Perylene-d12	1.000	1.000	0.0	95	0.01
87		Benzo(b)fluoranthene	1.394	1.173	15.9	75	0.01

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88	Benzo(k)fluoranthene	1.052	1.073	-2.0	119	0.01
89 C	Benzo(a)pyrene	1.128	1.077	4.5	89	0.01
90	Dibenzo(a,h)anthracene	1.047	0.933	10.9	84	0.02
91	Benzo(g,h,i)perylene	1.078	0.890	17.4	77	0.01

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