Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA F\Data\BF062717\
Data File : BF096217.D
Acq On : 27 Jun 2017 15:41
Operator : SJ/MA
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 28 15:59:04 2017
Quant Method : Z:\HPCHEM1\BNA F\METHODS\8270-BF062717.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 28 15:41:49 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	98	0.00
2		1,4-Dioxane	0.662	0.676	-2.1	101	0.00
3		Pyridine	1.816	1.923	-5.9	102	-0.02
4		n-Nitrosodimethylamine	0.898	0.930	-3.6	100	0.02
5	S	2-Fluorophenol	1.388	1.384	0.3	99	0.01
6		Aniline	2.318	2.306	0.5	97	0.01
7	S	Phenol-d6	1.681	1.686	-0.3	100	0.02
8		2-Chlorophenol	1.362	1.379	-1.2	99	0.00
9		Benzaldehyde	1.014	1.057	-4.2	100	0.00
10	C	Phenol	1.848	1.907	-3.2	99	0.02
11		bis(2-Chloroethyl)ether	1.512	1.513	-0.1	98	0.01
12		1,3-Dichlorobenzene	1.501	1.500	0.1	99	0.00
13	C	1,4-Dichlorobenzene	1.507	1.497	0.7	98	0.00
14		1,2-Dichlorobenzene	1.373	1.367	0.4	99	0.00
15		Benzyl Alcohol	1.134	1.175	-3.6	100	0.01
16		2,2'-oxybis(1-Chloropropane	3.194	3.200	-0.2	98	0.00
17		2-Methylphenol	1.181	1.205	-2.0	101	0.01
18		Hexachloroethane	0.544	0.572	-5.1	101	0.00
19	P	n-Nitroso-di-n-propylamine	1.007	1.029	-2.2	100	0.02
20		3+4-Methylphenols	1.347	1.347	0.0	100	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.416	0.409	1.7	99	-0.01
23	S	Nitrobenzene-d5	0.289	0.310	-7.3	103	-0.01
24		Nitrobenzene	0.320	0.353	-10.3	101	0.02
25		Isophorone	0.656	0.649	1.1	99	0.01
26	C	2-Nitrophenol	0.133	0.159	-19.5	111	-0.01
27		2,4-Dimethylphenol	0.296	0.292	1.4	98	0.00
28		bis(2-Chloroethoxy)methane	0.444	0.428	3.6	98	0.01
29	C	2,4-Dichlorophenol	0.253	0.258	-2.0	99	0.00
30		1,2,4-Trichlorobenzene	0.251	0.245	2.4	98	0.00
31		Naphthalene	0.964	0.932	3.3	99	0.00
32		Benzoic acid	0.182	0.214	-17.6	110	-0.05
33		4-Chloroaniline	0.399	0.397	0.5	99	0.00
34	C	Hexachlorobutadiene	0.124	0.121	2.4	96	0.00
35		Caprolactam	0.087	0.092	-5.7	101	0.06
36	C	4-Chloro-3-methylphenol	0.275	0.275	0.0	98	0.01
37		2-Methylnaphthalene	0.626	0.602	3.8	98	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
39		1,2,4,5-Tetrachlorobenzene	0.484	0.460	5.0	96	0.00
40	P	Hexachlorocyclopentadiene	0.217	0.237	-9.2	100	0.00
41	S	2,4,6-Tribromophenol	0.157	0.165	-5.1	100	0.00
42	C	2,4,6-Trichlorophenol	0.365	0.370	-1.4	98	0.00
43		2,4,5-Trichlorophenol	0.385	0.401	-4.2	97	0.01
44	S	2-Fluorobiphenyl	1.092	1.156	-5.9	98	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.535	1.539	-0.3	98	0.00
46		2-Chloronaphthalene	1.273	1.240	2.6	98	0.00
47		2-Nitroaniline	0.402	0.452	-12.4	107	-0.01
48		Acenaphthylene	2.112	2.062	2.4	99	0.00
49		Dimethylphthalate	1.431	1.409	1.5	98	0.01
50		2,6-Dinitrotoluene	0.298	0.323	-8.4	104	-0.02
51	C	Acenaphthene	1.243	1.205	3.1	99	0.01
52		3-Nitroaniline	0.381	0.410	-7.6	104	-0.01
53	P	2,4-Dinitrophenol	0.091	0.112	-23.1	118	-0.01
54		Dibenzofuran	1.623	1.611	0.7	97	0.00
55	P	4-Nitrophenol	0.308	0.336	-9.1	102	-0.02
56		2,4-Dinitrotoluene	0.375	0.417	-11.2	103	-0.02
57		Fluorene	1.188	1.158	2.5	98	0.00
58		2,3,4,6-Tetrachlorophenol	0.269	0.281	-4.5	99	0.00
59		Diethylphthalate	1.437	1.413	1.7	97	0.01
60		4-Chlorophenyl-phenylether	0.479	0.483	-0.8	99	0.00
61		4-Nitroaniline	0.332	0.349	-5.1	102	-0.02
62		Azobenzene	1.449	1.426	1.6	98	0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	99	0.00
64		4,6-Dinitro-2-methylphenol	0.091	0.110	-20.9	114	-0.01
65	c	n-Nitrosodiphenylamine	0.719	0.691	3.9	98	0.00
66		4-Bromophenyl-phenylether	0.206	0.201	2.4	98	0.00
67		Hexachlorobenzene	0.212	0.211	0.5	99	0.00
68		Atrazine	0.190	0.190	0.0	100	0.01
69	C	Pentachlorophenol	0.121	0.134	-10.7	101	0.00
70		Phenanthrene	1.080	1.073	0.6	99	-0.01
71		Anthracene	1.118	1.101	1.5	98	-0.01
72		Carbazole	1.234	1.169	5.3	97	0.01
73		Di-n-butylphthalate	1.303	1.288	1.2	97	0.00
74	C	Fluoranthene	1.111	1.080	2.8	99	0.01
75	I	Chrysene-d12	1.000	1.000	0.0	98	0.01
76		Benzidine	0.750	0.812	-8.3	100	0.00
77		Pyrene	1.644	1.653	-0.5	98	0.00
78	S	Terphenyl-d14	0.802	0.793	1.1	100	0.00
79		Butylbenzylphthalate	0.753	0.812	-7.8	100	0.00
80		Benzo(a)anthracene	1.219	1.241	-1.8	99	0.00
81		3,3'-Dichlorobenzidine	0.464	0.515	-11.0	101	0.00
82		Chrysene	1.261	1.306	-3.6	101	0.01
83		Bis(2-ethylhexyl)phthalate	0.819	0.891	-8.8	98	0.00
84	c	Di-n-octyl phthalate	1.686	1.845	-9.4	102	0.00
85		Indeno(1,2,3-cd)pyrene	1.047	1.083	-3.4	101	0.02
86	I	Perylene-d12	1.000	1.000	0.0	102	0.00
87		Benzo(b)fluoranthene	1.314	1.245	5.3	102	-0.02

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88	Benzo(k)fluoranthene	1.199	1.193	0.5	99	0.01
89 C	Benzo(a)pyrene	1.160	1.144	1.4	101	0.01
90	Dibenzo(a,h)anthracene	0.909	0.906	0.3	104	0.02
91	Benzo(g,h,i)perylene	0.944	0.901	4.6	102	0.02

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