Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\DATA\BF062415\
Data File : BF080129.D
Acq On : 24 Jun 2015 13:21
Operator : TP/IZ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 25 01:10:24 2015
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF061715.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 25 00:52:28 2015
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	84	-0.01
2		1,4-Dioxane	0.537	0.507	5.6	79	-0.01
3		Pyridine	1.428	1.341	6.1	79	0.00
4		n-Nitrosodimethylamine	0.679	0.574	15.5	67	-0.01
5	S	2-Fluorophenol	1.130	1.164	-3.0	84	-0.01
6		Aniline	2.068	2.065	0.1	80	0.00
7	S	Phenol-d6	1.503	1.540	-2.5	80	-0.01
8		2-Chlorophenol	1.306	1.306	0.0	81	-0.01
9		Benzaldehyde	0.868	0.716	17.5	67	-0.01
10	C	Phenol	1.641	1.794	-9.3	88	-0.01
11		bis(2-Chloroethyl)ether	1.302	1.310	-0.6	82	-0.01
12		1,3-Dichlorobenzene	1.569	1.595	-1.7	82	-0.01
13	C	1,4-Dichlorobenzene	1.492	1.495	-0.2	82	-0.01
14		1,2-Dichlorobenzene	1.452	1.503	-3.5	84	-0.01
15		Benzyl Alcohol	1.229	1.195	2.8	78	0.00
16		2,2'-oxybis(1-Chloropropane	1.932	1.866	3.4	82	-0.01
17		2-Methylphenol	1.115	1.130	-1.3	80	-0.01
18		Hexachloroethane	0.497	0.497	0.0	80	0.00
19	P	n-Nitroso-di-n-propylamine	1.044	0.991	5.1	83	-0.01
20		3+4-Methylphenols	1.440	1.419	1.5	79	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	84	-0.01
22		Acetophenone	0.466	0.457	1.9	77	-0.01
23	S	Nitrobenzene-d5	0.344	0.335	2.6	79	0.00
24		Nitrobenzene	0.355	0.357	-0.6	82	-0.01
25		Isophorone	0.691	0.674	2.5	79	-0.01
26	C	2-Nitrophenol	0.148	0.150	-1.4	82	-0.01
27		2,4-Dimethylphenol	0.309	0.290	6.1	80	-0.01
28		bis(2-Chloroethoxy)methane	0.406	0.402	1.0	80	0.00
29	C	2,4-Dichlorophenol	0.265	0.268	-1.1	81	-0.01
30		1,2,4-Trichlorobenzene	0.301	0.297	1.3	82	0.00
31		Naphthalene	1.046	1.038	0.8	80	0.00
32		Benzoic acid	0.187	0.123	34.2#	54	-0.01
33		4-Chloroaniline	0.448	0.413	7.8	79	-0.01
34	C	Hexachlorobutadiene	0.185	0.186	-0.5	87	-0.01
35		Caprolactam	0.086	0.083	3.5	75	-0.01
36	C	4-Chloro-3-methylphenol	0.299	0.295	1.3	78	-0.01
37		2-Methylnaphthalene	0.676	0.653	3.4	78	0.00
38	I	Acenaphthene-d10	1.000	1.000	0.0	85	-0.01
39		1,2,4,5-Tetrachlorobenzene	0.582	0.552	5.2	78	-0.01
40	P	Hexachlorocyclopentadiene	0.260	0.236	9.2	73	0.00
41	S	2,4,6-Tribromophenol	0.196	0.182	7.1	76	-0.01
42	C	2,4,6-Trichlorophenol	0.396	0.375	5.3	76	0.00
43		2,4,5-Trichlorophenol	0.456	0.450	1.3	79	-0.01
44	S	2-Fluorobiphenyl	1.402	1.335	4.8	79	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.627	1.526	6.2	78	-0.01
46		2-Chloronaphthalene	1.320	1.294	2.0	79	0.00
47		2-Nitroaniline	0.362	0.359	0.8	76	-0.01
48		Acenaphthylene	2.204	2.149	2.5	77	0.00
49		Dimethylphthalate	1.504	1.410	6.3	79	-0.01
50		2,6-Dinitrotoluene	0.301	0.314	-4.3	79	-0.01
51	C	Acenaphthene	1.348	1.326	1.6	80	-0.01
52		3-Nitroaniline	0.348	0.361	-3.7	81	-0.01
53	P	2,4-Dinitrophenol	0.106	0.106	0.0	79	-0.01
54		Dibenzofuran	1.913	1.816	5.1	77	-0.01
55	P	4-Nitrophenol	0.278	0.247	11.2	75	-0.01
56		2,4-Dinitrotoluene	0.383	0.384	-0.3	81	0.00
57		Fluorene	1.518	1.422	6.3	77	-0.01
58		2,3,4,6-Tetrachlorophenol	0.385	0.382	0.8	78	-0.01
59		Diethylphthalate	1.514	1.495	1.3	77	-0.01
60		4-Chlorophenyl-phenylether	0.729	0.691	5.2	79	-0.01
61		4-Nitroaniline	0.359	0.346	3.6	75	0.00
62		Azobenzene	1.541	1.495	3.0	76	-0.01
63	I	Phenanthrene-d10	1.000	1.000	0.0	85	-0.02
64		4,6-Dinitro-2-methylphenol	0.094	0.099	-5.3	82	-0.01
65	c	n-Nitrosodiphenylamine	0.651	0.620	4.8	76	0.00
66		4-Bromophenyl-phenylether	0.213	0.200	6.1	76	-0.01
67		Hexachlorobenzene	0.236	0.217	8.1	74	-0.02
68		Atrazine	0.206	0.182	11.7	69	-0.02
69	C	Pentachlorophenol	0.134	0.121	9.7	77	-0.01
70		Phenanthrene	1.211	1.092	9.8	75	-0.02
71		Anthracene	1.166	1.138	2.4	82	-0.02
72		Carbazole	1.113	1.004	9.8	74	-0.02
73		Di-n-butylphthalate	1.286	1.155	10.2	78	-0.02
74	C	Fluoranthene	1.290	1.223	5.2	81	-0.05
75	I	Chrysene-d12	1.000	1.000	0.0	81	-0.07
76		Benzidine	0.559	0.546	2.3	76	-0.05
77		Pyrene	1.231	1.211	1.6	78	-0.05
78	S	Terphenyl-d14	0.856	0.822	4.0	78	-0.06
79		Butylbenzylphthalate	0.495	0.487	1.6	75	-0.07
80		Benzo(a)anthracene	1.218	1.191	2.2	79	-0.07
81		3,3'-Dichlorobenzidine	0.420	0.402	4.3	76	-0.07
82		Chrysene	1.124	1.090	3.0	77	-0.07
83		Bis(2-ethylhexyl)phthalate	0.782	0.757	3.2	75	-0.07
84	c	Di-n-octyl phthalate	1.339	1.289	3.7	75	-0.09
85		Indeno(1,2,3-cd)pyrene	1.417	1.411	0.4	80	-0.10
86	I	Perylene-d12	1.000	1.000	0.0	82	-0.09
87		Benzo(b)fluoranthene	1.211	1.174	3.1	77	-0.09

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88	Benzo(k)fluoranthene	1.233	1.209	1.9	80	-0.09
89 C	Benzo(a)pyrene	1.150	1.098	4.5	77	-0.09
90	Dibenzo(a,h)anthracene	1.177	1.148	2.5	79	-0.10
91	Benzo(g,h,i)perylene	1.188	1.135	4.5	77	-0.10

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