Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA_F\Data\BF060116\
Data File : BF087642.D
Acq On : 2 Jun 2016 6:31
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_F
LabSampleId : SSTDCCC040

Quant Time: Jun 02 12:28:11 2016
Quant Method : Z:\HPCHEM1\BNA_F\METHODS\8270-BF052316.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 25 16:26:52 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	120	-0.09
2		1,4-Dioxane	0.601	0.487	19.0	99	-0.10
3		Pyridine	1.313	1.119	14.8	95	-0.13
4		n-Nitrosodimethylamine	1.012	0.905	10.6	104	-0.10
5	S	2-Fluorophenol	1.138	1.068	6.2	106	-0.09
6		Aniline	1.990	1.791	10.0	103	-0.09
7	S	Phenol-d6	1.538	1.333	13.3	103	-0.07
8		2-Chlorophenol	1.419	1.230	13.3	103	-0.08
9		Benzaldehyde	0.849	0.740	12.8	114	-0.08
10	C	Phenol	1.679	1.479	11.9	114	-0.06
11		bis(2-Chloroethyl)ether	1.359	1.130	16.9	101	-0.09
12		1,3-Dichlorobenzene	1.548	1.294	16.4	101	-0.09
13	C	1,4-Dichlorobenzene	1.589	1.352	14.9	103	-0.10
14		1,2-Dichlorobenzene	1.470	1.254	14.7	105	-0.10
15		Benzyl Alcohol	1.121	1.096	2.2	109	-0.07
16		2,2'-oxybis(1-Chloropropane	3.058	2.386	22.0	95	-0.09
17		2-Methylphenol	1.138	1.002	12.0	106	-0.07
18		Hexachloroethane	0.593	0.512	13.7	104	-0.10
19	P	n-Nitroso-di-n-propylamine	1.147	0.939	18.1	99	-0.08
20		3+4-Methylphenols	1.375	1.236	10.1	102	-0.07
21	I	Naphthalene-d8	1.000	1.000	0.0	116	-0.09
22		Acetophenone	0.478	0.429	10.3	103	-0.08
23	S	Nitrobenzene-d5	0.397	0.358	9.8	101	-0.08
24		Nitrobenzene	0.419	0.377	10.0	100	-0.09
25		Isophorone	0.712	0.636	10.7	103	-0.09
26	C	2-Nitrophenol	0.171	0.156	8.8	98	-0.09
27		2,4-Dimethylphenol	0.297	0.252	15.2	96	-0.08
28		bis(2-Chloroethoxy)methane	0.401	0.340	15.2	100	-0.09
29	C	2,4-Dichlorophenol	0.268	0.240	10.4	100	-0.07
30		1,2,4-Trichlorobenzene	0.307	0.263	14.3	99	-0.10
31		Naphthalene	1.002	0.882	12.0	105	-0.09
32		Benzoic acid	0.141	0.135	4.3	102	0.00
33		4-Chloroaniline	0.369	0.345	6.5	105	-0.09
34	C	Hexachlorobutadiene	0.186	0.161	13.4	100	-0.11
35		Caprolactam	0.078	0.076	2.6	109	-0.05
36	C	4-Chloro-3-methylphenol	0.303	0.286	5.6	104	-0.06
37		2-Methylnaphthalene	0.626	0.546	12.8	102	-0.10
38	I	Acenaphthene-d10	1.000	1.000	0.0	123	-0.10
39		1,2,4,5-Tetrachlorobenzene	0.640	0.524	18.1	100	-0.10
40	P	Hexachlorocyclopentadiene	0.206	0.097	52.9#	51	-0.10
41	S	2,4,6-Tribromophenol	0.192	0.164	14.6	101	-0.09
42	C	2,4,6-Trichlorophenol	0.370	0.308	16.8	98	-0.08
43		2,4,5-Trichlorophenol	0.376	0.287	23.7	92	-0.07
44	S	2-Fluorobiphenyl	1.419	1.143	19.5	101	-0.10

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.684	1.415	16.0	107	-0.10
46		2-Chloronaphthalene	1.288	1.072	16.8	103	-0.09
47		2-Nitroaniline	0.464	0.437	5.8	111	-0.08
48		Acenaphthylene	2.040	1.594	21.9	97	-0.10
49		Dimethylphthalate	1.602	1.308	18.4	100	-0.09
50		2,6-Dinitrotoluene	0.323	0.277	14.2	105	-0.08
51	C	Acenaphthene	1.254	1.034	17.5	105	-0.09
52		3-Nitroaniline	0.331	0.302	8.8	110	-0.07
53	P	2,4-Dinitrophenol	0.138	0.099	28.3#	76	-0.03
54		Dibenzofuran	1.697	1.410	16.9	105	-0.09
55	P	4-Nitrophenol	0.158	0.092	41.8#	70	-0.01
56		2,4-Dinitrotoluene	0.431	0.394	8.6	107	-0.07
57		Fluorene	1.472	1.178	20.0	100	-0.10
58		2,3,4,6-Tetrachlorophenol	0.289	0.218	24.6	87	-0.08
59		Diethylphthalate	1.558	1.320	15.3	106	-0.09
60		4-Chlorophenyl-phenylether	0.718	0.569	20.8	99	-0.10
61		4-Nitroaniline	0.309	0.280	9.4	100	-0.06
62		Azobenzene	1.616	1.438	11.0	104	-0.10
63	I	Phenanthrene-d10	1.000	1.000	0.0	122	-0.09
64		4,6-Dinitro-2-methylphenol	0.127	0.101	20.5	87	-0.06
65	c	n-Nitrosodiphenylamine	0.647	0.553	14.5	102	-0.09
66		4-Bromophenyl-phenylether	0.219	0.189	13.7	104	-0.10
67		Hexachlorobenzene	0.229	0.199	13.1	104	-0.10
68		Atrazine	0.206	0.080	61.2#	46#	-0.09
69	C	Pentachlorophenol	0.061	0.032	47.5#	56	-0.07
70		Phenanthrene	1.108	0.955	13.8	106	-0.09
71		Anthracene	1.119	0.952	14.9	105	-0.09
72		Carbazole	0.955	0.815	14.7	103	-0.08
73		Di-n-butylphthalate	1.234	1.059	14.2	99	-0.09
74	C	Fluoranthene	1.111	0.965	13.1	104	-0.08
75	I	Chrysene-d12	1.000	1.000	0.0	114	-0.08
76		Benzidine	0.515	0.399	22.5	80	-0.08
77		Pyrene	1.440	1.347	6.5	106	-0.08
78	S	Terphenyl-d14	0.941	0.906	3.7	108	-0.08
79		Butylbenzylphthalate	0.638	0.627	1.7	104	-0.08
80		Benzo(a)anthracene	1.138	0.965	15.2	96	-0.07
81		3,3'-Dichlorobenzidine	0.628	0.577	8.1	108	-0.08
82		Chrysene	1.129	0.929	17.7	96	-0.08
83		Bis(2-ethylhexyl)phthalate	0.932	0.759	18.6	91	-0.09
84	c	Di-n-octyl phthalate	1.421	1.166	17.9	87	-0.09
85		Indeno(1,2,3-cd)pyrene	0.905	0.879	2.9	109	-0.08
86	I	Perylene-d12	1.000	1.000	0.0	116	-0.06
87		Benzo(b)fluoranthene	1.274	0.887	30.4#	75	-0.06

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(k)fluoranthene	1.097	1.085	1.1	135	-0.07
89 C	Benzo(a)pyrene	1.102	0.926	16.0	100	-0.06
90	Dibenzo(a,h)anthracene	0.940	0.882	6.2	107	-0.08
91	Benzo(g,h,i)perylene	0.927	0.851	8.2	105	-0.08

(#) = Out of Range SPCC's out = 0 CCC's out = 1