Evaluate Continuing Calibration Report

Data Path : Z:\HPCHEM1\BNA M\DATA\BM122515\
Data File : BM003798.D
Acq On : 25 Dec 2015 03:15
Operator : UM/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Dec 25 06:52:20 2015
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM120915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Dec 09 18:52:47 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	66	-0.04
2		1,4-Dioxane	0.467	0.457	2.1	62	-0.04
3		Pyridine	1.308	1.313	-0.4	63	-0.04
4		n-Nitrosodimethylamine	0.419	0.460	-9.8	68	-0.04
5	S	2-Fluorophenol	1.125	1.115	0.9	62	-0.03
6		Aniline	2.056	2.095	-1.9	65	-0.04
7	S	Phenol-d6	1.562	1.591	-1.9	64	-0.04
8		2-Chlorophenol	1.424	1.412	0.8	63	-0.04
9		Benzaldehyde	0.766	0.784	-2.3	61	-0.04
10	C	Phenol	1.657	1.701	-2.7	65	-0.03
11		bis(2-Chloroethyl)ether	1.341	1.339	0.1	63	-0.04
12		1,3-Dichlorobenzene	1.565	1.526	2.5	62	-0.04
13	C	1,4-Dichlorobenzene	1.613	1.586	1.7	63	-0.04
14		1,2-Dichlorobenzene	1.541	1.509	2.1	63	-0.04
15		Benzyl Alcohol	1.081	1.163	-7.6	66	-0.04
16		2,2'-oxybis(1-Chloropropane	1.433	1.561	-8.9	70	-0.04
17		2-Methylphenol	1.159	1.212	-4.6	66	-0.04
18		Hexachloroethane	0.554	0.555	-0.2	63	-0.05
19	P	n-Nitroso-di-n-propylamine	1.025	1.110	-8.3	68	-0.04
20		3+4-Methylphenols	1.603	1.664	-3.8	65	-0.04
21	I	Naphthalene-d8	1.000	1.000	0.0	69	-0.05
22		Acetophenone	0.490	0.465	5.1	63	-0.04
23	S	Nitrobenzene-d5	0.323	0.330	-2.2	67	-0.04
24		Nitrobenzene	0.349	0.352	-0.9	67	-0.04
25		Isophorone	0.652	0.669	-2.6	67	-0.04
26	C	2-Nitrophenol	0.176	0.168	4.5	64	-0.04
27		2,4-Dimethylphenol	0.307	0.300	2.3	65	-0.04
28		bis(2-Chloroethoxy)methane	0.392	0.379	3.3	65	-0.04
29	C	2,4-Dichlorophenol	0.293	0.288	1.7	64	-0.04
30		1,2,4-Trichlorobenzene	0.327	0.307	6.1	63	-0.04
31		Naphthalene	1.047	1.006	3.9	65	-0.04
32		Benzoic acid	0.196	0.174	11.2	63	-0.02
33		4-Chloroaniline	0.432	0.427	1.2	66	-0.04
34	C	Hexachlorobutadiene	0.194	0.183	5.7	63	-0.05
35		Caprolactam	0.097	0.106	-9.3	76	-0.04
36	C	4-Chloro-3-methylphenol	0.330	0.340	-3.0	68	-0.04
37		2-Methylnaphthalene	0.718	0.697	2.9	66	-0.04
38	I	Acenaphthene-d10	1.000	1.000	0.0	74	-0.04
39		1,2,4,5-Tetrachlorobenzene	0.643	0.570	11.4	63	-0.04
40	P	Hexachlorocyclopentadiene	0.359	0.282	21.4	56	-0.04
41	S	2,4,6-Tribromophenol	0.200	0.209	-4.5	75	-0.04
42	C	2,4,6-Trichlorophenol	0.419	0.407	2.9	67	-0.04
43		2,4,5-Trichlorophenol	0.441	0.432	2.0	69	-0.04
44	S	2-Fluorobiphenyl	1.468	1.317	10.3	64	-0.04

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.774	1.596	10.0	65	-0.04
46		2-Chloronaphthalene	1.303	1.219	6.4	68	-0.04
47		2-Nitroaniline	0.323	0.341	-5.6	76	-0.03
48		Acenaphthylene	2.172	2.105	3.1	69	-0.04
49		Dimethylphthalate	1.648	1.609	2.4	71	-0.04
50		2,6-Dinitrotoluene	0.344	0.350	-1.7	72	-0.04
51	C	Acenaphthene	1.259	1.196	5.0	69	-0.04
52		3-Nitroaniline	0.352	0.380	-8.0	76	-0.03
53	P	2,4-Dinitrophenol	0.165	0.147	10.9	66	-0.02
54		Dibenzofuran	1.818	1.756	3.4	71	-0.04
55	P	4-Nitrophenol	0.291	0.304	-4.5	78	-0.02
56		2,4-Dinitrotoluene	0.446	0.486	-9.0	77	-0.04
57		Fluorene	1.507	1.490	1.1	73	-0.04
58		2,3,4,6-Tetrachlorophenol	0.339	0.349	-2.9	73	-0.04
59		Diethylphthalate	1.618	1.643	-1.5	75	-0.04
60		4-Chlorophenyl-phenylether	0.762	0.722	5.2	71	-0.04
61		4-Nitroaniline	0.359	0.408	-13.6	85	-0.03
62		Azobenzene	1.269	1.380	-8.7	78	-0.04
63	I	Phenanthrene-d10	1.000	1.000	0.0	85	-0.04
64		4,6-Dinitro-2-methylphenol	0.121	0.112	7.4	77	-0.03
65	c	n-Nitrosodiphenylamine	0.632	0.586	7.3	74	-0.04
66		4-Bromophenyl-phenylether	0.216	0.199	7.9	74	-0.04
67		Hexachlorobenzene	0.232	0.216	6.9	76	-0.04
68		Atrazine	0.197	0.207	-5.1	83	-0.04
69	C	Pentachlorophenol	0.132	0.125	5.3	77	-0.04
70		Phenanthrene	1.122	1.078	3.9	80	-0.04
71		Anthracene	1.115	1.098	1.5	81	-0.04
72		Carbazole	0.972	1.026	-5.6	87	-0.04
73		Di-n-butylphthalate	1.136	1.255	-10.5	89	-0.04
74	C	Fluoranthene	1.157	1.269	-9.7	93	-0.04
75	I	Chrysene-d12	1.000	1.000	0.0	110	-0.04
76		Benzidine	0.641	0.636	0.8	100	-0.04
77		Pyrene	1.403	1.193	15.0	93	-0.04
78	S	Terphenyl-d14	0.848	0.763	10.0	98	-0.04
79		Butylbenzylphthalate	0.540	0.543	-0.6	109	-0.04
80		Benzo(a)anthracene	1.199	1.139	5.0	101	-0.04
81		3,3'-Dichlorobenzidine	0.382	0.415	-8.6	109	-0.04
82		Chrysene	1.154	1.089	5.6	103	-0.04
83		Bis(2-ethylhexyl)phthalate	0.741	0.772	-4.2	111	-0.03
84	c	Di-n-octyl phthalate	1.224	1.398	-14.2	121	-0.05
85		Indeno(1,2,3-cd)pyrene	1.213	1.062	12.4	89	-0.08
86	I	Perylene-d12	1.000	1.000	0.0	110	-0.05
87		Benzo(b)fluoranthene	1.216	1.220	-0.3	109	-0.05

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88	Benzo(k)fluoranthene	1.187	1.125	5.2	102	-0.05
89 C	Benzo(a)pyrene	1.144	1.102	3.7	102	-0.05
90	Dibenzo(a,h)anthracene	1.117	0.942	15.7	89	-0.09
91	Benzo(g,h,i)perylene	1.128	0.903	19.9	84	-0.08

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