Evaluate Continuing Calibration Report

Data Path : \\74.0.250.170\svoasrv\HPCHEM1\BNA M\Data\BM090717\
Data File : BM011455.D
Acq On : 07 Sep 2017 15:25
Operator : SJ/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Sep 08 12:18:54 2017
Quant Method : Z:\HPCHEM1\BNA M\METHODS\8270-BM083117.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Aug 31 18:39:59 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	85	-0.01
2		1,4-Dioxane	0.495	0.482	2.6	83	-0.01
3		Pyridine	1.517	1.422	6.3	78	0.00
4		n-Nitrosodimethylamine	0.774	0.813	-5.0	89	0.00
5	S	2-Fluorophenol	1.188	1.149	3.3	82	-0.01
6		Aniline	2.238	2.069	7.6	77	-0.01
7	S	Phenol-d6	1.648	1.682	-2.1	85	0.00
8		2-Chlorophenol	1.401	1.376	1.8	82	-0.02
9		Benzaldehyde	0.958	0.812	15.2	76	-0.01
10	C	Phenol	1.811	1.807	0.2	83	-0.01
11		bis(2-Chloroethyl)ether	1.448	1.466	-1.2	85	-0.01
12		1,3-Dichlorobenzene	1.595	1.549	2.9	82	-0.01
13	C	1,4-Dichlorobenzene	1.622	1.579	2.7	82	-0.01
14		1,2-Dichlorobenzene	1.570	1.540	1.9	83	-0.02
15		Benzyl Alcohol	1.191	1.021	14.3	71	-0.01
16		2,2'-oxybis(1-Chloropropane	2.421	2.470	-2.0	86	-0.02
17		2-Methylphenol	1.160	1.164	-0.3	84	-0.02
18		Hexachloroethane	0.555	0.543	2.2	82	-0.02
19	P	n-Nitroso-di-n-propylamine	1.210	1.241	-2.6	86	-0.02
20		3+4-Methylphenols	1.609	1.602	0.4	83	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	87	-0.02
22		Acetophenone	0.497	0.479	3.6	84	-0.01
23	S	Nitrobenzene-d5	0.303	0.333	-9.9	92	-0.01
24		Nitrobenzene	0.334	0.356	-6.6	89	-0.02
25		Isophorone	0.680	0.640	5.9	81	-0.02
26	C	2-Nitrophenol	0.133	0.146	-9.8	96	-0.01
27		2,4-Dimethylphenol	0.317	0.299	5.7	82	-0.02
28		bis(2-Chloroethoxy)methane	0.464	0.451	2.8	85	-0.02
29	C	2,4-Dichlorophenol	0.297	0.289	2.7	83	-0.02
30		1,2,4-Trichlorobenzene	0.328	0.312	4.9	83	-0.02
31		Naphthalene	1.080	1.043	3.4	85	-0.02
32		Benzoic acid	0.196	0.207	-5.6	91	-0.01
33		4-Chloroaniline	0.472	0.411	12.9	75	-0.01
34	C	Hexachlorobutadiene	0.187	0.175	6.4	82	-0.02
35		Caprolactam	0.107	0.096	10.3	78	0.00
36	C	4-Chloro-3-methylphenol	0.348	0.338	2.9	84	-0.01
37		2-Methylnaphthalene	0.752	0.732	2.7	85	-0.02
38	I	Acenaphthene-d10	1.000	1.000	0.0	88	-0.02
39		1,2,4,5-Tetrachlorobenzene	0.602	0.588	2.3	85	-0.02
40	P	Hexachlorocyclopentadiene	0.276	0.276	0.0	88	-0.02
41	S	2,4,6-Tribromophenol	0.231	0.224	3.0	83	-0.01
42	C	2,4,6-Trichlorophenol	0.398	0.395	0.8	85	-0.01
43		2,4,5-Trichlorophenol	0.405	0.400	1.2	84	-0.01
44	S	2-Fluorobiphenyl	1.387	1.425	-2.7	89	-0.02

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45		1,1'-Biphenyl	1.776	1.745	1.7	87	-0.02
46		2-Chloronaphthalene	1.319	1.280	3.0	85	-0.02
47		2-Nitroaniline	0.387	0.445	-15.0	100	-0.01
48		Acenaphthylene	2.055	1.997	2.8	85	-0.02
49		Dimethylphthalate	1.589	1.496	5.9	83	-0.02
50		2,6-Dinitrotoluene	0.301	0.302	-0.3	87	-0.02
51	C	Acenaphthene	1.335	1.281	4.0	85	-0.02
52		3-Nitroaniline	0.391	0.384	1.8	85	-0.01
53	P	2,4-Dinitrophenol	0.097	0.148	-52.6#	140	-0.01
54		Dibenzofuran	1.840	1.789	2.8	86	-0.01
55	P	4-Nitrophenol	0.304	0.286	5.9	82	0.00
56		2,4-Dinitrotoluene	0.394	0.409	-3.8	90	-0.01
57		Fluorene	1.604	1.637	-2.1	90	-0.02
58		2,3,4,6-Tetrachlorophenol	0.327	0.329	-0.6	86	-0.02
59		Diethylphthalate	1.637	1.531	6.5	83	-0.02
60		4-Chlorophenyl-phenylether	0.748	0.761	-1.7	90	-0.02
61		4-Nitroaniline	0.406	0.399	1.7	85	-0.01
62		Azobenzene	1.476	1.497	-1.4	89	-0.02
63	I	Phenanthrene-d10	1.000	1.000	0.0	89	-0.01
64		4,6-Dinitro-2-methylphenol	0.072	0.097	-34.7#	122	-0.01
65	c	n-Nitrosodiphenylamine	0.630	0.617	2.1	87	-0.01
66		4-Bromophenyl-phenylether	0.210	0.200	4.8	84	-0.02
67		Hexachlorobenzene	0.247	0.229	7.3	82	-0.02
68		Atrazine	0.185	0.159	14.1	76	-0.02
69	C	Pentachlorophenol	0.136	0.139	-2.2	87	-0.02
70		Phenanthrene	1.114	1.114	0.0	88	-0.02
71		Anthracene	1.120	1.121	-0.1	88	-0.01
72		Carbazole	1.079	1.056	2.1	86	-0.01
73		Di-n-butylphthalate	1.254	1.242	1.0	84	-0.02
74	C	Fluoranthene	1.271	1.301	-2.4	89	-0.02
75	I	Chrysene-d12	1.000	1.000	0.0	96	-0.02
76		Benzidine	0.395	0.271	31.4#	68	-0.01
77		Pyrene	1.226	1.182	3.6	90	-0.01
78	S	Terphenyl-d14	0.806	0.801	0.6	94	-0.02
79		Butylbenzylphthalate	0.541	0.498	7.9	85	-0.02
80		Benzo(a)anthracene	1.119	1.108	1.0	95	-0.02
81		3,3'-Dichlorobenzidine	0.425	0.409	3.8	88	-0.02
82		Chrysene	1.054	1.029	2.4	93	-0.02
83		Bis(2-ethylhexyl)phthalate	0.794	0.774	2.5	90	-0.03
84	c	Di-n-octyl phthalate	1.338	1.336	0.1	88	-0.04
85		Indeno(1,2,3-cd)pyrene	0.983	1.031	-4.9	101	-0.04
86	I	Perylene-d12	1.000	1.000	0.0	98	-0.02
87		Benzo(b)fluoranthene	1.228	1.190	3.1	94	-0.02

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88	Benzo(k)fluoranthene	1.177	1.177	0.0	95	-0.02
89 C	Benzo(a)pyrene	1.107	1.093	1.3	94	-0.02
90	Dibenzo(a,h)anthracene	0.948	0.971	-2.4	99	-0.04
91	Benzo(g,h,i)perylene	0.916	0.924	-0.9	98	-0.04

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