Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM120218\
Data File : BM017871.D
Acq On : 02 Dec 2018 00:30
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Dec 03 01:19:13 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM111218.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Nov 12 18:37:25 2018
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	135	-0.04
2		1,4-Dioxane	0.494	0.415	16.0	116	-0.03
3		Pyridine	1.448	1.276	11.9	117	-0.02
4		n-Nitrosodimethylamine	0.638	0.550	13.8	114	-0.02
5	S	2-Fluorophenol	1.185	1.163	1.9	131	-0.03
6		Aniline	2.053	2.070	-0.8	132	-0.03
7	S	Phenol-d6	1.655	1.689	-2.1	134	-0.02
8		2-Chlorophenol	1.418	1.482	-4.5	140	-0.03
9		Benzaldehyde	1.206	0.996	17.4	109	-0.04
10	C	Phenol	1.737	1.765	-1.6	135	-0.03
11		bis(2-Chloroethyl)ether	1.371	1.429	-4.2	138	-0.03
12		1,3-Dichlorobenzene	1.591	1.656	-4.1	141	-0.04
13	C	1,4-Dichlorobenzene	1.631	1.695	-3.9	140	-0.04
14		1,2-Dichlorobenzene	1.581	1.657	-4.8	141	-0.04
15		Benzyl Alcohol	1.318	1.380	-4.7	135	-0.03
16		2,2'-oxybis(1-Chloropropane	2.088	1.630	21.9	104	-0.03
17		2-Methylphenol	1.170	1.221	-4.4	137	-0.03
18		Hexachloroethane	0.575	0.638	-11.0	145	-0.04
19	P	n-Nitroso-di-n-propylamine	1.188	1.257	-5.8	137	-0.03
20		3+4-Methylphenols	1.626	1.709	-5.1	136	-0.02
21	I	Naphthalene-d8	1.000	1.000	0.0	135	-0.04
22		Acetophenone	0.499	0.520	-4.2	137	-0.03
23	S	Nitrobenzene-d5	0.387	0.397	-2.6	135	-0.03
24		Nitrobenzene	0.393	0.395	-0.5	135	-0.03
25		Isophorone	0.598	0.627	-4.8	137	-0.03
26	C	2-Nitrophenol	0.181	0.196	-8.3	140	-0.04
27		2,4-Dimethylphenol	0.261	0.271	-3.8	136	-0.03
28		bis(2-Chloroethoxy)methane	0.405	0.424	-4.7	137	-0.03
29	C	2,4-Dichlorophenol	0.303	0.327	-7.9	139	-0.04
30		1,2,4-Trichlorobenzene	0.352	0.371	-5.4	141	-0.03
31		Naphthalene	0.983	1.019	-3.7	138	-0.04
32		Benzoic acid	0.236	0.246	-4.2	137	0.00
33		4-Chloroaniline	0.431	0.446	-3.5	133	-0.04
34	C	Hexachlorobutadiene	0.219	0.236	-7.8	145	-0.04
35		Caprolactam	0.112	0.111	0.9	133	0.00
36	C	4-Chloro-3-methylphenol	0.350	0.364	-4.0	136	-0.03
37		2-Methylnaphthalene	0.695	0.725	-4.3	138	-0.04
38		1-Methylnaphthalene	0.681	0.703	-3.2	137	-0.04
39	I	Acenaphthene-d10	1.000	1.000	0.0	126	-0.03
40		1,2,4,5-Tetrachlorobenzene	0.705	0.776	-10.1	138	-0.04
41	P	Hexachlorocyclopentadiene	0.364	0.379	-4.1	129	-0.04
42	S	2,4,6-Tribromophenol	0.287	0.294	-2.4	125	-0.02
43	C	2,4,6-Trichlorophenol	0.442	0.496	-12.2	137	-0.03
44		2,4,5-Trichlorophenol	0.466	0.509	-9.2	134	-0.04

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.542	1.674	-8.6	136	-0.03
46		1,1'-Biphenyl	1.795	1.937	-7.9	136	-0.04
47		2-Chloronaphthalene	1.332	1.454	-9.2	137	-0.03
48		2-Nitroaniline	0.412	0.439	-6.6	128	-0.03
49		Acenaphthylene	2.002	2.124	-6.1	133	-0.03
50		Dimethylphthalate	1.626	1.710	-5.2	132	-0.02
51		2,6-Dinitrotoluene	0.362	0.385	-6.4	131	-0.02
52	C	Acenaphthene	1.229	1.288	-4.8	132	-0.03
53		3-Nitroaniline	0.394	0.402	-2.0	128	-0.02
54	P	2,4-Dinitrophenol	0.186	0.276	-48.4#	180#	-0.02
55		Dibenzofuran	2.008	2.070	-3.1	131	-0.03
56	P	4-Nitrophenol	0.294	0.326	-10.9	137	-0.02
57		2,4-Dinitrotoluene	0.488	0.527	-8.0	129	-0.02
58		Fluorene	1.537	1.599	-4.0	131	-0.03
59		2,3,4,6-Tetrachlorophenol	0.429	0.444	-3.5	127	-0.03
60		Diethylphthalate	1.757	1.818	-3.5	134	-0.02
61		4-Chlorophenyl-phenylether	0.873	0.914	-4.7	132	-0.03
62		4-Nitroaniline	0.372	0.375	-0.8	125	-0.02
63		Azobenzene	1.624	1.707	-5.1	130	-0.03
64	I	Phenanthrene-d10	1.000	1.000	0.0	113	-0.03
65		4,6-Dinitro-2-methylphenol	0.139	0.157	-12.9	123	-0.02
66	c	n-Nitrosodiphenylamine	0.643	0.747	-16.2	128	-0.03
67		4-Bromophenyl-phenylether	0.237	0.273	-15.2	127	-0.03
68		Hexachlorobenzene	0.267	0.290	-8.6	122	-0.02
69		Atrazine	0.237	0.253	-6.8	115	-0.02
70	C	Pentachlorophenol	0.187	0.210	-12.3	124	-0.03
71		Phenanthrene	1.085	1.136	-4.7	117	-0.02
72		Anthracene	1.056	1.126	-6.6	117	-0.03
73		Carbazole	1.067	1.103	-3.4	112	-0.02
74		Di-n-butylphthalate	1.188	1.382	-16.3	124	-0.02
75	C	Fluoranthene	1.228	1.110	9.6	99	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	84	-0.02
77		Benzidine	0.651	0.674	-3.5	82	-0.02
78		Pyrene	1.234	1.428	-15.7	96	-0.02
79	S	Terphenyl-d14	0.882	1.040	-17.9	98	-0.02
80		Butylbenzylphthalate	0.501	0.619	-23.6	97	-0.03
81		Benzo(a)anthracene	1.152	1.220	-5.9	87	-0.02
82		3,3'-Dichlorobenzidine	0.445	0.505	-13.5	89	-0.02
83		Chrysene	1.119	1.168	-4.4	86	-0.02
84		Bis(2-ethylhexyl)phthalate	0.741	0.880	-18.8	97	-0.02
85	c	Di-n-octyl phthalate	1.186	1.347	-13.6	92	-0.02
86		Indeno(1,2,3-cd)pyrene	0.894	1.369	-53.1#	123	-0.05
87	I	Perylene-d12	1.000	1.000	0.0	96	-0.04

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88	Benzo(b)fluoranthene	1.171	1.165	0.5	95	-0.03
89	Benzo(k)fluoranthene	1.184	1.161	1.9	91	-0.03
90 C	Benzo(a)pyrene	1.068	1.130	-5.8	100	-0.04
91	Dibenzo(a,h)anthracene	0.898	1.172	-30.5#	124	-0.05
92	Benzo(g,h,i)perylene	0.844	1.134	-34.4#	128	-0.05

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