Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM030119\
Data File : BM018985.D
Acq On : 01 Mar 2019 12:56
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Mar 01 14:18:55 2019
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM021119.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Mar 01 14:12:30 2019
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	97	-0.04
2		1,4-Dioxane	0.532	0.505	5.1	92	-0.02
3		Pyridine	1.450	1.540	-6.2	94	-0.02
4		n-Nitrosodimethylamine	0.369	0.402	-8.9	101	-0.02
5	S	2-Fluorophenol	1.221	1.248	-2.2	97	-0.03
6		Aniline	2.107	2.272	-7.8	102	-0.03
7	S	Phenol-d6	1.720	1.879	-9.2	103	-0.02
8		2-Chlorophenol	1.340	1.429	-6.6	102	-0.03
9		Benzaldehyde	0.941	0.937	0.4	97	-0.03
10	C	Phenol	1.809	1.935	-7.0	102	-0.03
11		bis(2-Chloroethyl)ether	1.380	1.484	-7.5	102	-0.03
12		1,3-Dichlorobenzene	1.507	1.560	-3.5	100	-0.03
13	C	1,4-Dichlorobenzene	1.534	1.592	-3.8	101	-0.03
14		1,2-Dichlorobenzene	1.474	1.553	-5.4	101	-0.03
15		Benzyl Alcohol	1.366	1.530	-12.0	104	-0.03
16		2,2'-oxybis(1-Chloropropane	1.336	1.481	-10.9	108	-0.03
17		2-Methylphenol	1.152	1.245	-8.1	102	-0.03
18		Hexachloroethane	0.560	0.595	-6.2	102	-0.04
19	P	n-Nitroso-di-n-propylamine	1.331	1.500	-12.7	105	-0.03
20		3+4-Methylphenols	1.590	1.758	-10.6	103	-0.03
21	I	Naphthalene-d8	1.000	1.000	0.0	99	-0.04
22		Acetophenone	0.549	0.579	-5.5	104	-0.03
23	S	Nitrobenzene-d5	0.433	0.457	-5.5	103	-0.03
24		Nitrobenzene	0.449	0.464	-3.3	101	-0.03
25		Isophorone	0.690	0.757	-9.7	106	-0.03
26	C	2-Nitrophenol	0.179	0.200	-11.7	106	-0.03
27		2,4-Dimethylphenol	0.261	0.283	-8.4	106	-0.03
28		bis(2-Chloroethoxy)methane	0.446	0.477	-7.0	105	-0.03
29	C	2,4-Dichlorophenol	0.300	0.323	-7.7	103	-0.04
30		1,2,4-Trichlorobenzene	0.346	0.359	-3.8	103	-0.03
31		Naphthalene	0.975	1.009	-3.5	104	-0.04
32		Benzoic acid	0.228	0.226	0.9	98	0.00
33		4-Chloroaniline	0.436	0.473	-8.5	105	-0.03
34	C	Hexachlorobutadiene	0.201	0.201	0.0	101	-0.04
35		Caprolactam	0.122	0.137	-12.3	108	-0.02
36	C	4-Chloro-3-methylphenol	0.362	0.391	-8.0	104	-0.02
37		2-Methylnaphthalene	0.687	0.727	-5.8	105	-0.03
38		1-Methylnaphthalene	0.672	0.709	-5.5	105	-0.03
39	I	Acenaphthene-d10	1.000	1.000	0.0	99	-0.02
40		1,2,4,5-Tetrachlorobenzene	0.640	0.661	-3.3	103	-0.03
41	P	Hexachlorocyclopentadiene	0.327	0.325	0.6	98	-0.03
42	S	2,4,6-Tribromophenol	0.288	0.327	-13.5	110	-0.02
43	C	2,4,6-Trichlorophenol	0.424	0.457	-7.8	104	-0.03
44		2,4,5-Trichlorophenol	0.444	0.476	-7.2	103	-0.03

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.507	1.594	-5.8	106	-0.03
46		1,1'-Biphenyl	1.722	1.808	-5.0	106	-0.03
47		2-Chloronaphthalene	1.330	1.396	-5.0	105	-0.03
48		2-Nitroaniline	0.442	0.503	-13.8	108	-0.02
49		Acenaphthylene	1.980	2.115	-6.8	105	-0.03
50		Dimethylphthalate	1.667	1.772	-6.3	106	-0.02
51		2,6-Dinitrotoluene	0.361	0.401	-11.1	107	-0.02
52	C	Acenaphthene	1.214	1.282	-5.6	105	-0.02
53		3-Nitroaniline	0.408	0.435	-6.6	105	-0.02
54	P	2,4-Dinitrophenol	0.200	0.222	-11.0	110	-0.02
55		Dibenzofuran	1.995	2.091	-4.8	105	-0.03
56	P	4-Nitrophenol	0.326	0.404	-23.9	122	-0.02
57		2,4-Dinitrotoluene	0.497	0.557	-12.1	106	-0.02
58		Fluorene	1.527	1.639	-7.3	108	-0.02
59		2,3,4,6-Tetrachlorophenol	0.392	0.409	-4.3	101	-0.02
60		Diethylphthalate	1.719	1.829	-6.4	106	-0.02
61		4-Chlorophenyl-phenylether	0.856	0.888	-3.7	104	-0.02
62		4-Nitroaniline	0.400	0.424	-6.0	104	-0.02
63		Azobenzene	1.844	1.998	-8.4	106	-0.02
64	I	Phenanthrene-d10	1.000	1.000	0.0	99	-0.02
65		4,6-Dinitro-2-methylphenol	0.140	0.153	-9.3	107	-0.02
66	c	n-Nitrosodiphenylamine	0.632	0.683	-8.1	106	-0.02
67		4-Bromophenyl-phenylether	0.224	0.237	-5.8	104	-0.03
68		Hexachlorobenzene	0.260	0.274	-5.4	105	-0.02
69		Atrazine	0.204	0.183	10.3	88	-0.02
70	C	Pentachlorophenol	0.168	0.179	-6.5	100	-0.02
71		Phenanthrene	1.075	1.126	-4.7	105	-0.02
72		Anthracene	1.046	1.123	-7.4	106	-0.02
73		Carbazole	1.093	1.156	-5.8	104	-0.02
74		Di-n-butylphthalate	1.257	1.392	-10.7	107	-0.02
75	C	Fluoranthene	1.242	1.264	-1.8	100	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	86	-0.02
77		Benzidine	0.450	0.514	-14.2	80	-0.02
78		Pyrene	1.162	1.353	-16.4	100	-0.02
79	S	Terphenyl-d14	0.970	1.148	-18.4	99	-0.02
80		Butylbenzylphthalate	0.529	0.642	-21.4	99	-0.02
81		Benzo(a)anthracene	1.166	1.215	-4.2	89	-0.02
82		3,3'-Dichlorobenzidine	0.461	0.489	-6.1	88	-0.02
83		Chrysene	1.117	1.156	-3.5	89	-0.02
84		Bis(2-ethylhexyl)phthalate	0.775	0.893	-15.2	96	-0.02
85	c	Di-n-octyl phthalate	1.302	1.372	-5.4	86	-0.02
86		Indeno(1,2,3-cd)pyrene	1.290	1.089	15.6	66	-0.05
87	I	Perylene-d12	1.000	1.000	0.0	70	-0.04

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.144	1.232	-7.7	76	-0.03
89	Benzo(k)fluoranthene	1.139	1.220	-7.1	76	-0.03
90 C	Benzo(a)pyrene	1.084	1.136	-4.8	73	-0.04
91	Dibenzo(a,h)anthracene	1.074	1.077	-0.3	67	-0.05
92	Benzo(g,h,i)perylene	1.000	1.010	-1.0	66	-0.05

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