Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM120418\
Data File : BM017881.D
Acq On : 03 Dec 2018 19:28
Operator : JU/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Dec 04 07:17:46 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	86	-0.04
2		1,4-Dioxane	0.494	0.445	9.9	79	-0.03
3		Pyridine	1.448	1.359	6.1	79	-0.03
4		n-Nitrosodimethylamine	0.638	0.494	22.6	65	-0.02
5	S	2-Fluorophenol	1.185	1.187	-0.2	85	-0.03
6		Aniline	2.053	2.062	-0.4	84	-0.04
7	S	Phenol-d6	1.655	1.674	-1.1	84	-0.04
8		2-Chlorophenol	1.418	1.470	-3.7	88	-0.04
9		Benzaldehyde	1.206	0.994	17.6	69	-0.04
10	C	Phenol	1.737	1.737	0.0	84	-0.04
11		bis(2-Chloroethyl)ether	1.371	1.375	-0.3	84	-0.04
12		1,3-Dichlorobenzene	1.591	1.662	-4.5	90	-0.04
13	C	1,4-Dichlorobenzene	1.631	1.688	-3.5	88	-0.04
14		1,2-Dichlorobenzene	1.581	1.622	-2.6	87	-0.04
15		Benzyl Alcohol	1.318	1.310	0.6	81	-0.04
16		2,2'-oxybis(1-Chloropropane	2.088	1.567	25.0	63	-0.04
17		2-Methylphenol	1.170	1.184	-1.2	84	-0.04
18		Hexachloroethane	0.575	0.602	-4.7	87	-0.04
19	P	n-Nitroso-di-n-propylamine	1.188	1.190	-0.2	82	-0.04
20		3+4-Methylphenols	1.626	1.680	-3.3	85	-0.04
21	I	Naphthalene-d8	1.000	1.000	0.0	85	-0.04
22		Acetophenone	0.499	0.506	-1.4	84	-0.04
23	S	Nitrobenzene-d5	0.387	0.381	1.6	81	-0.04
24		Nitrobenzene	0.393	0.378	3.8	81	-0.04
25		Isophorone	0.598	0.617	-3.2	85	-0.04
26	C	2-Nitrophenol	0.181	0.196	-8.3	88	-0.04
27		2,4-Dimethylphenol	0.261	0.272	-4.2	86	-0.04
28		bis(2-Chloroethoxy)methane	0.405	0.411	-1.5	83	-0.04
29	C	2,4-Dichlorophenol	0.303	0.320	-5.6	85	-0.04
30		1,2,4-Trichlorobenzene	0.352	0.368	-4.5	87	-0.04
31		Naphthalene	0.983	1.002	-1.9	85	-0.04
32		Benzoic acid	0.236	0.219	7.2	76	-0.03
33		4-Chloroaniline	0.431	0.449	-4.2	84	-0.04
34	C	Hexachlorobutadiene	0.219	0.227	-3.7	88	-0.04
35		Caprolactam	0.112	0.115	-2.7	87	-0.04
36	C	4-Chloro-3-methylphenol	0.350	0.356	-1.7	83	-0.04
37		2-Methylnaphthalene	0.695	0.720	-3.6	86	-0.04
38		1-Methylnaphthalene	0.681	0.706	-3.7	86	-0.04
39	I	Acenaphthene-d10	1.000	1.000	0.0	83	-0.04
40		1,2,4,5-Tetrachlorobenzene	0.705	0.725	-2.8	85	-0.04
41	P	Hexachlorocyclopentadiene	0.364	0.357	1.9	80	-0.04
42	S	2,4,6-Tribromophenol	0.287	0.312	-8.7	87	-0.03
43	C	2,4,6-Trichlorophenol	0.442	0.468	-5.9	85	-0.04
44		2,4,5-Trichlorophenol	0.466	0.489	-4.9	85	-0.04

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.542	1.614	-4.7	86	-0.04
46		1,1'-Biphenyl	1.795	1.846	-2.8	85	-0.04
47		2-Chloronaphthalene	1.332	1.377	-3.4	85	-0.04
48		2-Nitroaniline	0.412	0.387	6.1	74	-0.04
49		Acenaphthylene	2.002	2.099	-4.8	86	-0.04
50		Dimethylphthalate	1.626	1.653	-1.7	84	-0.03
51		2,6-Dinitrotoluene	0.362	0.381	-5.2	85	-0.03
52	C	Acenaphthene	1.229	1.294	-5.3	87	-0.04
53		3-Nitroaniline	0.394	0.405	-2.8	85	-0.03
54	P	2,4-Dinitrophenol	0.186	0.199	-7.0	85	-0.03
55		Dibenzofuran	2.008	2.088	-4.0	87	-0.03
56	P	4-Nitrophenol	0.294	0.315	-7.1	87	-0.03
57		2,4-Dinitrotoluene	0.488	0.532	-9.0	85	-0.03
58		Fluorene	1.537	1.606	-4.5	86	-0.03
59		2,3,4,6-Tetrachlorophenol	0.429	0.456	-6.3	85	-0.04
60		Diethylphthalate	1.757	1.795	-2.2	87	-0.03
61		4-Chlorophenyl-phenylether	0.873	0.908	-4.0	86	-0.03
62		4-Nitroaniline	0.372	0.393	-5.6	86	-0.02
63		Azobenzene	1.624	1.666	-2.6	83	-0.03
64	I	Phenanthrene-d10	1.000	1.000	0.0	83	-0.03
65		4,6-Dinitro-2-methylphenol	0.139	0.150	-7.9	86	-0.03
66	c	n-Nitrosodiphenylamine	0.643	0.686	-6.7	86	-0.03
67		4-Bromophenyl-phenylether	0.237	0.251	-5.9	85	-0.04
68		Hexachlorobenzene	0.267	0.281	-5.2	86	-0.03
69		Atrazine	0.237	0.244	-3.0	81	-0.03
70	C	Pentachlorophenol	0.187	0.201	-7.5	87	-0.03
71		Phenanthrene	1.085	1.124	-3.6	85	-0.03
72		Anthracene	1.056	1.119	-6.0	85	-0.03
73		Carbazole	1.067	1.116	-4.6	83	-0.03
74		Di-n-butylphthalate	1.188	1.342	-13.0	89	-0.03
75	C	Fluoranthene	1.228	1.254	-2.1	82	-0.03
76	I	Chrysene-d12	1.000	1.000	0.0	73	-0.02
77		Benzidine	0.651	0.697	-7.1	73	-0.02
78		Pyrene	1.234	1.396	-13.1	82	-0.03
79	S	Terphenyl-d14	0.882	1.004	-13.8	82	-0.02
80		Butylbenzylphthalate	0.501	0.632	-26.1#	86	-0.03
81		Benzo(a)anthracene	1.152	1.201	-4.3	74	-0.02
82		3,3'-Dichlorobenzidine	0.445	0.498	-11.9	76	-0.03
83		Chrysene	1.119	1.150	-2.8	74	-0.03
84		Bis(2-ethylhexyl)phthalate	0.741	0.915	-23.5	87	-0.02
85	c	Di-n-octyl phthalate	1.186	1.367	-15.3	81	-0.03
86		Indeno(1,2,3-cd)pyrene	0.894	1.049	-17.3	82	-0.05
87	I	Perylene-d12	1.000	1.000	0.0	70	-0.04

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88	Benzo(b)fluoranthene	1.171	1.213	-3.6	72	-0.04
89	Benzo(k)fluoranthene	1.184	1.199	-1.3	68	-0.03
90 C	Benzo(a)pyrene	1.068	1.116	-4.5	71	-0.04
91	Dibenzo(a,h)anthracene	0.898	1.087	-21.0	84	-0.05
92	Benzo(g,h,i)perylene	0.844	1.034	-22.5	85	-0.05

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