Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA M\Data\BM102218\
Data File : BM017274.D
Acq On : 22 Oct 2018 12:05
Operator : MJ/SJ
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_M
LabSampleId : SSTDCCC040

Quant Time: Oct 22 13:05:14 2018
Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\8270-BM101718.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Oct 17 14:23:17 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	118	-0.01
2		1,4-Dioxane	0.604	0.573	5.1	114	-0.01
3		Pyridine	1.389	1.467	-5.6	118	-0.01
4		n-Nitrosodimethylamine	0.558	0.564	-1.1	120	0.00
5	S	2-Fluorophenol	1.182	1.259	-6.5	122	-0.01
6		Aniline	2.014	2.206	-9.5	124	-0.01
7	S	Phenol-d6	1.610	1.774	-10.2	124	-0.01
8		2-Chlorophenol	1.368	1.519	-11.0	126	-0.02
9		Benzaldehyde	1.027	1.045	-1.8	121	-0.02
10	C	Phenol	1.733	1.872	-8.0	122	-0.01
11		bis(2-Chloroethyl)ether	1.382	1.487	-7.6	124	-0.02
12		1,3-Dichlorobenzene	1.595	1.680	-5.3	123	-0.01
13	C	1,4-Dichlorobenzene	1.629	1.726	-6.0	124	-0.01
14		1,2-Dichlorobenzene	1.570	1.647	-4.9	122	-0.01
15		Benzyl Alcohol	1.177	1.436	-22.0	134	-0.02
16		2,2'-oxybis(1-Chloropropane	1.930	1.974	-2.3	119	-0.02
17		2-Methylphenol	1.113	1.264	-13.6	128	-0.02
18		Hexachloroethane	0.558	0.602	-7.9	124	-0.02
19	P	n-Nitroso-di-n-propylamine	1.060	1.269	-19.7	131	-0.01
20		3+4-Methylphenols	1.522	1.808	-18.8	132	-0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	125	-0.01
22		Acetophenone	0.507	0.528	-4.1	128	-0.02
23	S	Nitrobenzene-d5	0.342	0.383	-12.0	130	-0.02
24		Nitrobenzene	0.362	0.389	-7.5	126	-0.02
25		Isophorone	0.565	0.638	-12.9	133	-0.02
26	C	2-Nitrophenol	0.156	0.204	-30.8#	159#	-0.01
27		2,4-Dimethylphenol	0.250	0.281	-12.4	134	-0.01
28		bis(2-Chloroethoxy)methane	0.397	0.429	-8.1	130	-0.02
29	C	2,4-Dichlorophenol	0.290	0.336	-15.9	138	-0.01
30		1,2,4-Trichlorobenzene	0.348	0.366	-5.2	132	-0.02
31		Naphthalene	0.980	1.016	-3.7	130	-0.02
32		Benzoic acid	0.186	0.227	-22.0	155#	0.00
33		4-Chloroaniline	0.423	0.480	-13.5	136	-0.02
34	C	Hexachlorobutadiene	0.220	0.227	-3.2	130	-0.02
35		Caprolactam	0.106	0.126	-18.9	145	0.00
36	C	4-Chloro-3-methylphenol	0.327	0.383	-17.1	139	-0.01
37		2-Methylnaphthalene	0.671	0.747	-11.3	135	-0.02
38		1-Methylnaphthalene	0.653	0.726	-11.2	135	-0.02
39	I	Acenaphthene-d10	1.000	1.000	0.0	132	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.704	0.733	-4.1	137	-0.02
41	P	Hexachlorocyclopentadiene	0.340	0.392	-15.3	144	-0.02
42	S	2,4,6-Tribromophenol	0.270	0.312	-15.6	151#	-0.01
43	C	2,4,6-Trichlorophenol	0.409	0.482	-17.8	144	-0.02
44		2,4,5-Trichlorophenol	0.438	0.503	-14.8	142	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
45	S	2-Fluorobiphenyl	1.503	1.562	-3.9	136	-0.01
46		1,1'-Biphenyl	1.800	1.872	-4.0	137	-0.01
47		2-Chloronaphthalene	1.324	1.370	-3.5	134	-0.02
48		2-Nitroaniline	0.342	0.414	-21.1	150#	-0.01
49		Acenaphthylene	1.919	2.080	-8.4	136	-0.01
50		Dimethylphthalate	1.543	1.706	-10.6	141	-0.01
51		2,6-Dinitrotoluene	0.340	0.388	-14.1	145	-0.01
52	C	Acenaphthene	1.205	1.268	-5.2	136	-0.02
53		3-Nitroaniline	0.368	0.419	-13.9	145	-0.01
54	P	2,4-Dinitrophenol	0.162	0.197	-21.6	155#	-0.01
55		Dibenzofuran	2.004	2.079	-3.7	136	-0.02
56	P	4-Nitrophenol	0.299	0.318	-6.4	134	-0.01
57		2,4-Dinitrotoluene	0.456	0.536	-17.5	147	-0.01
58		Fluorene	1.483	1.582	-6.7	138	-0.02
59		2,3,4,6-Tetrachlorophenol	0.400	0.462	-15.5	143	-0.02
60		Diethylphthalate	1.530	1.740	-13.7	142	-0.01
61		4-Chlorophenyl-phenylether	0.846	0.904	-6.9	138	-0.01
62		4-Nitroaniline	0.381	0.433	-13.6	143	-0.01
63		Azobenzene	1.487	1.598	-7.5	133	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	131	-0.01
65		4,6-Dinitro-2-methylphenol	0.116	0.129	-11.2	141	-0.01
66	c	n-Nitrosodiphenylamine	0.605	0.673	-11.2	140	-0.01
67		4-Bromophenyl-phenylether	0.220	0.251	-14.1	144	-0.01
68		Hexachlorobenzene	0.255	0.278	-9.0	142	-0.02
69		Atrazine	0.217	0.249	-14.7	146	-0.01
70	C	Pentachlorophenol	0.175	0.195	-11.4	141	-0.02
71		Phenanthrene	1.079	1.126	-4.4	136	-0.01
72		Anthracene	1.025	1.128	-10.0	138	-0.02
73		Carbazole	1.047	1.136	-8.5	138	-0.01
74		Di-n-butylphthalate	1.105	1.315	-19.0	151#	-0.01
75	C	Fluoranthene	1.214	1.319	-8.6	137	-0.02
76	I	Chrysene-d12	1.000	1.000	0.0	125	-0.01
77		Benzidine	0.507	0.633	-24.9	142	-0.01
78		Pyrene	1.118	1.282	-14.7	136	-0.01
79	S	Terphenyl-d14	0.887	1.007	-13.5	137	0.00
80		Butylbenzylphthalate	0.384	0.561	-46.1#	179#	-0.01
81		Benzo(a)anthracene	1.125	1.210	-7.6	131	-0.01
82		3,3'-Dichlorobenzidine	0.419	0.471	-12.4	135	-0.01
83		Chrysene	1.114	1.148	-3.1	128	-0.01
84		Bis(2-ethylhexyl)phthalate	0.601	0.809	-34.6#	161#	-0.01
85	c	Di-n-octyl phthalate	1.018	1.303	-28.0#	155#	-0.01
86		Indeno(1,2,3-cd)pyrene	1.246	1.001	19.7	100	-0.02
87	I	Perylene-d12	1.000	1.000	0.0	110	-0.02

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.093	1.213	-11.0	120	-0.01
89	Benzo(k)fluoranthene	1.097	1.205	-9.8	118	-0.01
90 C	Benzo(a)pyrene	1.038	1.108	-6.7	115	-0.02
91	Dibenzo(a,h)anthracene	1.031	0.950	7.9	100	-0.03
92	Benzo(g,h,i)perylene	1.022	0.905	11.4	97	-0.03

(#) = Out of Range SPCC's out = 0 CCC's out = 2