Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP080221\
Data File : BP006462.D
Acq On : 02 Aug 2021 21:47
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 16 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Aug 03 05:30:07 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP080221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 03 05:26:21 2021
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	104	0.00
2		1,4-Dioxane	0.568	0.559	1.6	103	0.00
3		Pyridine	1.641	1.675	-2.1	105	0.00
4		n-Nitrosodimethylamine	0.620	0.638	-2.9	106	0.00
5	S	2-Fluorophenol	1.282	1.291	-0.7	104	0.00
6		Aniline	1.979	2.078	-5.0	106	0.00
7	S	Phenol-d6	1.811	1.901	-5.0	107	0.00
8		2-Chlorophenol	1.387	1.431	-3.2	107	0.00
9		Benzaldehyde	1.135	1.100	3.1	107	0.00
10	C	Phenol	1.824	1.895	-3.9	106	0.00
11		bis(2-Chloroethyl)ether	1.396	1.437	-2.9	107	0.00
12		1,3-Dichlorobenzene	1.464	1.462	0.1	105	0.00
13	C	1,4-Dichlorobenzene	1.481	1.478	0.2	104	0.00
14		1,2-Dichlorobenzene	1.423	1.428	-0.4	106	0.00
15		Benzyl Alcohol	1.356	1.452	-7.1	109	0.00
16		2,2'-oxybis(1-Chloropropane	1.648	1.704	-3.4	108	0.00
17		2-Methylphenol	1.141	1.222	-7.1	107	0.00
18		Hexachloroethane	0.580	0.584	-0.7	106	0.00
19	P	n-Nitroso-di-n-propylamine	1.201	1.314	-9.4	110	0.00
20		3+4-Methylphenols	1.557	1.680	-7.9	109	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	108	0.00
22		Acetophenone	0.527	0.526	0.2	108	0.00
23	S	Nitrobenzene-d5	0.424	0.429	-1.2	109	0.00
24		Nitrobenzene	0.440	0.447	-1.6	109	0.00
25		Isophorone	0.757	0.797	-5.3	110	0.00
26	C	2-Nitrophenol	0.158	0.166	-5.1	110	0.00
27		2,4-Dimethylphenol	0.270	0.281	-4.1	108	0.00
28		bis(2-Chloroethoxy)methane	0.414	0.416	-0.5	109	0.00
29	C	2,4-Dichlorophenol	0.273	0.282	-3.3	109	0.00
30		1,2,4-Trichlorobenzene	0.296	0.291	1.7	109	0.00
31		Naphthalene	1.065	1.062	0.3	109	0.00
32		Benzoic acid	0.205	0.215	-4.9	113	0.00
33		4-Chloroaniline	0.426	0.443	-4.0	108	0.00
34	C	Hexachlorobutadiene	0.173	0.168	2.9	108	0.00
35		Caprolactam	0.100	0.111	-11.0	110	0.00
36	C	4-Chloro-3-methylphenol	0.340	0.372	-9.4	111	0.00
37		2-Methylnaphthalene	0.715	0.740	-3.5	110	0.00
38		1-Methylnaphthalene	0.679	0.701	-3.2	110	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
40		1,2,4,5-Tetrachlorobenzene	0.521	0.495	5.0	110	0.00
41	P	Hexachlorocyclopentadiene	0.276	0.270	2.2	110	0.00
42	S	2,4,6-Tribromophenol	0.205	0.212	-3.4	110	0.00
43	C	2,4,6-Trichlorophenol	0.350	0.354	-1.1	111	0.00
44		2,4,5-Trichlorophenol	0.387	0.392	-1.3	111	0.00
45	S	2-Fluorobiphenyl	1.334	1.279	4.1	110	0.00
46		1,1'-Biphenyl	1.512	1.453	3.9	110	0.00
47		2-Chloronaphthalene	1.164	1.119	3.9	110	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.390	0.421	-7.9	112	0.00
49		Acenaphthylene	1.850	1.858	-0.4	110	0.00
50		Dimethylphthalate	1.438	1.444	-0.4	110	0.00
51		2,6-Dinitrotoluene	0.313	0.328	-4.8	110	0.00
52	C	Acenaphthene	1.141	1.119	1.9	110	0.00
53		3-Nitroaniline	0.347	0.378	-8.9	111	0.00
54	P	2,4-Dinitrophenol	0.170	0.170	0.0	108	0.00
55		Dibenzofuran	1.787	1.765	1.2	110	0.00
56	P	4-Nitrophenol	0.315	0.336	-6.7	110	0.00
57		2,4-Dinitrotoluene	0.418	0.455	-8.9	111	0.00
58		Fluorene	1.419	1.422	-0.2	110	0.00
59		2,3,4,6-Tetrachlorophenol	0.325	0.333	-2.5	110	0.00
60		Diethylphthalate	1.495	1.535	-2.7	110	0.00
61		4-Chlorophenyl-phenylether	0.645	0.637	1.2	110	0.00
62		4-Nitroaniline	0.367	0.410	-11.7	110	0.00
63		Azobenzene	1.726	1.814	-5.1	111	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	110	0.00
65		4,6-Dinitro-2-methylphenol	0.120	0.118	1.7	109	0.00
66	c	n-Nitrosodiphenylamine	0.619	0.606	2.1	110	0.00
67		4-Bromophenyl-phenylether	0.192	0.183	4.7	108	0.00
68		Hexachlorobenzene	0.216	0.211	2.3	112	0.00
69		Atrazine	0.191	0.197	-3.1	109	0.00
70	C	Pentachlorophenol	0.134	0.141	-5.2	112	0.00
71		Phenanthrene	1.122	1.098	2.1	109	0.00
72		Anthracene	1.075	1.074	0.1	109	0.00
73		Carbazole	1.092	1.109	-1.6	109	0.00
74		Di-n-butylphthalate	1.233	1.279	-3.7	109	0.00
75	C	Fluoranthene	1.248	1.269	-1.7	108	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	107	0.00
77		Benzidine	0.492	0.529	-7.5	102	0.00
78		Pyrene	1.319	1.304	1.1	108	0.00
79	S	Terphenyl-d14	0.993	0.963	3.0	108	0.00
80		Butylbenzylphthalate	0.570	0.600	-5.3	109	0.00
81		Benzo(a)anthracene	1.288	1.286	0.2	108	0.00
82		3,3'-Dichlorobenzidine	0.341	0.348	-2.1	106	0.00
83		Chrysene	1.255	1.242	1.0	107	0.00
84		Bis(2-ethylhexyl)phthalate	0.880	0.885	-0.6	106	0.00
85	c	Di-n-octyl phthalate	1.490	1.480	0.7	107	0.00
86	I	Perylene-d12	1.000	1.000	0.0	107	0.00
87		Indeno(1,2,3-cd)pyrene	1.435	1.449	-1.0	107	0.00
88		Benzo(b)fluoranthene	1.284	1.288	-0.3	109	0.00
89		Benzo(k)fluoranthene	1.245	1.240	0.4	106	0.00
90	C	Benzo(a)pyrene	1.149	1.177	-2.4	108	0.00
91		Dibenzo(a,h)anthracene	1.243	1.251	-0.6	107	0.00
92		Benzo(g,h,i)perylene	1.191	1.205	-1.2	107	0.00

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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