Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011922\
Data File : BP008846.D
Acq On : 19 Jan 2022 10:44
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jan 19 13:06:01 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 14 18:05:00 2022
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	91	0.00
2		1,4-Dioxane	0.523	0.522	0.2	108	0.00
3		Pyridine	1.096	1.147	-4.7	106	0.00
4		n-Nitrosodimethylamine	0.514	0.543	-5.6	107	0.00
5	S	2-Fluorophenol	1.293	1.314	-1.6	105	0.00
6		Aniline	1.955	1.978	-1.2	102	0.00
7	S	Phenol-d6	1.566	1.601	-2.2	103	0.00
8		2-Chlorophenol	1.377	1.404	-2.0	104	0.00
9		Benzaldehyde	1.047	1.001	4.4	98	0.00
10	C	Phenol	1.597	1.623	-1.6	103	0.00
11		bis(2-Chloroethyl)ether	1.270	1.295	-2.0	104	0.00
12		1,3-Dichlorobenzene	1.640	1.647	-0.4	103	0.00
13	C	1,4-Dichlorobenzene	1.662	1.669	-0.4	104	0.00
14		1,2-Dichlorobenzene	1.585	1.585	0.0	103	0.00
15		Benzyl Alcohol	1.099	1.115	-1.5	101	0.00
16		2,2'-oxybis(1-Chloropropane	1.421	1.458	-2.6	103	0.00
17		2-Methylphenol	1.070	1.077	-0.7	101	0.00
18		Hexachloroethane	0.564	0.576	-2.1	104	0.00
19	P	n-Nitroso-di-n-propylamine	0.915	0.937	-2.4	100	0.00
20		3+4-Methylphenols	1.420	1.433	-0.9	100	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	89	0.00
22		Acetophenone	0.509	0.519	-2.0	100	0.00
23	S	Nitrobenzene-d5	0.352	0.367	-4.3	102	0.00
24		Nitrobenzene	0.356	0.372	-4.5	102	0.00
25		Isophorone	0.637	0.657	-3.1	100	0.00
26	C	2-Nitrophenol	0.187	0.194	-3.7	100	0.00
27		2,4-Dimethylphenol	0.319	0.327	-2.5	99	0.00
28		bis(2-Chloroethoxy)methane	0.415	0.424	-2.2	100	0.00
29	C	2,4-Dichlorophenol	0.348	0.357	-2.6	100	0.00
30		1,2,4-Trichlorobenzene	0.408	0.412	-1.0	101	0.00
31		Naphthalene	1.086	1.090	-0.4	100	0.00
32		Benzoic acid	0.183	0.188	-2.7	94	-0.01
33		4-Chloroaniline	0.443	0.448	-1.1	98	0.00
34	C	Hexachlorobutadiene	0.255	0.256	-0.4	100	0.00
35		Caprolactam	0.096	0.098	-2.1	98	0.00
36	C	4-Chloro-3-methylphenol	0.314	0.320	-1.9	98	0.00
37		2-Methylnaphthalene	0.794	0.810	-2.0	100	0.00
38		1-Methylnaphthalene	0.729	0.740	-1.5	99	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	88	0.00
40		1,2,4,5-Tetrachlorobenzene	0.729	0.734	-0.7	100	0.00
41	P	Hexachlorocyclopentadiene	0.373	0.381	-2.1	93	0.00
42	S	2,4,6-Tribromophenol	0.291	0.298	-2.4	100	0.00
43	C	2,4,6-Trichlorophenol	0.451	0.458	-1.6	98	0.00
44		2,4,5-Trichlorophenol	0.485	0.495	-2.1	99	0.00
45	S	2-Fluorobiphenyl	1.517	1.518	-0.1	100	0.00
46		1,1'-Biphenyl	1.627	1.635	-0.5	100	0.00
47		2-Chloronaphthalene	1.255	1.270	-1.2	100	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.279	0.289	-3.6	100	0.00
49		Acenaphthylene	1.896	1.922	-1.4	100	0.00
50		Dimethylphthalate	1.485	1.500	-1.0	100	0.00
51		2,6-Dinitrotoluene	0.315	0.329	-4.4	101	0.00
52	C	Acenaphthene	1.124	1.145	-1.9	101	0.00
53		3-Nitroaniline	0.310	0.321	-3.5	100	0.00
54	P	2,4-Dinitrophenol	0.127	0.141	-11.0	104	0.00
55		Dibenzofuran	1.832	1.856	-1.3	100	0.00
56	P	4-Nitrophenol	0.223	0.236	-5.8	100	0.00
57		2,4-Dinitrotoluene	0.405	0.431	-6.4	101	0.00
58		Fluorene	1.447	1.480	-2.3	101	0.00
59		2,3,4,6-Tetrachlorophenol	0.400	0.412	-3.0	100	0.00
60		Diethylphthalate	1.409	1.462	-3.8	102	0.00
61		4-Chlorophenyl-phenylether	0.788	0.811	-2.9	101	0.00
62		4-Nitroaniline	0.300	0.317	-5.7	102	0.00
63		Azobenzene	1.170	1.238	-5.8	102	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	92	0.00
65		4,6-Dinitro-2-methylphenol	0.100	0.112	-12.0	107	0.00
66	c	n-Nitrosodiphenylamine	0.644	0.648	-0.6	102	0.00
67		4-Bromophenyl-phenylether	0.250	0.252	-0.8	101	0.00
68		Hexachlorobenzene	0.286	0.282	1.4	101	0.00
69		Atrazine	0.239	0.247	-3.3	105	0.00
70	C	Pentachlorophenol	0.152	0.158	-3.9	104	0.00
71		Phenanthrene	1.126	1.129	-0.3	103	0.00
72		Anthracene	1.130	1.164	-3.0	105	0.00
73		Carbazole	0.971	1.004	-3.4	107	0.00
74		Di-n-butylphthalate	1.169	1.241	-6.2	107	0.00
75	C	Fluoranthene	1.252	1.307	-4.4	110	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	104	0.00
77		Benzidine	0.707	0.767	-8.5	112	0.00
78		Pyrene	1.396	1.391	0.4	111	0.00
79	S	Terphenyl-d14	1.185	1.157	2.4	109	0.00
80		Butylbenzylphthalate	0.561	0.557	0.7	112	0.00
81		Benzo(a)anthracene	1.346	1.355	-0.7	117	0.00
82		3,3'-Dichlorobenzidine	0.581	0.603	-3.8	119	0.00
83		Chrysene	1.304	1.309	-0.4	117	0.00
84		Bis(2-ethylhexyl)phthalate	0.869	0.879	-1.2	113	0.00
85	c	Di-n-octyl phthalate	1.470	1.482	-0.8	124	0.00
86	I	Perylene-d12	1.000	1.000	0.0	108	0.00
87		Indeno(1,2,3-cd)pyrene	1.648	1.706	-3.5	126	-0.01
88		Benzo(b)fluoranthene	1.336	1.385	-3.7	127	0.00
89		Benzo(k)fluoranthene	1.293	1.322	-2.2	128	0.00
90	C	Benzo(a)pyrene	1.289	1.347	-4.5	128	0.00
91		Dibenzo(a,h)anthracene	1.401	1.445	-3.1	125	-0.01
92		Benzo(g,h,i)perylene	1.368	1.393	-1.8	123	-0.01

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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