Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA P\Data\BP020720\
Data File : BP001723.D
Acq On : 07 Feb 2020 16:23
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Feb 08 01:19:47 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP020320.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 03 15:16:18 2020
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	100	0.00
2		1,4-Dioxane	0.462	0.464	-0.4	101	0.00
3		Pyridine	1.319	1.376	-4.3	100	0.00
4		n-Nitrosodimethylamine	0.503	0.498	1.0	97	0.00
5	S	2-Fluorophenol	1.124	1.163	-3.5	101	0.00
6		Aniline	1.803	1.831	-1.6	97	0.00
7	S	Phenol-d6	1.466	1.516	-3.4	99	0.00
8		2-Chlorophenol	1.232	1.282	-4.1	100	0.00
9		Benzaldehyde	0.795	0.753	5.3	99	0.00
10	C	Phenol	1.494	1.533	-2.6	98	0.00
11		bis(2-Chloroethyl)ether	1.200	1.219	-1.6	98	0.00
12		1,3-Dichlorobenzene	1.502	1.516	-0.9	99	0.00
13	C	1,4-Dichlorobenzene	1.510	1.552	-2.8	101	0.00
14		1,2-Dichlorobenzene	1.435	1.457	-1.5	99	0.00
15		Benzyl Alcohol	0.993	1.029	-3.6	99	0.00
16		2,2'-oxybis(1-Chloropropane	1.498	1.501	-0.2	96	-0.01
17		2-Methylphenol	0.994	1.021	-2.7	98	0.00
18		Hexachloroethane	0.525	0.542	-3.2	102	0.00
19	P	n-Nitroso-di-n-propylamine	0.891	0.898	-0.8	96	0.00
20		3+4-Methylphenols	1.326	1.369	-3.2	99	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	98	0.00
22		Acetophenone	0.490	0.490	0.0	97	0.00
23	S	Nitrobenzene-d5	0.327	0.336	-2.8	99	0.00
24		Nitrobenzene	0.333	0.337	-1.2	98	0.00
25		Isophorone	0.612	0.611	0.2	96	0.00
26	C	2-Nitrophenol	0.115	0.124	-7.8	113	0.00
27		2,4-Dimethylphenol	0.243	0.245	-0.8	99	0.00
28		bis(2-Chloroethoxy)methane	0.416	0.414	0.5	97	0.00
29	C	2,4-Dichlorophenol	0.283	0.293	-3.5	100	0.00
30		1,2,4-Trichlorobenzene	0.347	0.348	-0.3	99	0.00
31		Naphthalene	1.020	1.023	-0.3	98	0.00
32		Benzoic acid	0.117	0.127	-8.5	122	0.00
33		4-Chloroaniline	0.435	0.433	0.5	96	0.00
34	C	Hexachlorobutadiene	0.206	0.210	-1.9	101	0.00
35		Caprolactam	0.086	0.091	-5.8	103	0.00
36	C	4-Chloro-3-methylphenol	0.289	0.290	-0.3	97	0.00
37		2-Methylnaphthalene	0.714	0.719	-0.7	98	0.00
38		1-Methylnaphthalene	0.677	0.676	0.1	98	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	98	0.00
40		1,2,4,5-Tetrachlorobenzene	0.636	0.631	0.8	97	0.00
41	P	Hexachlorocyclopentadiene	0.275	0.285	-3.6	103	0.00
42	S	2,4,6-Tribromophenol	0.205	0.223	-8.8	107	0.00
43	C	2,4,6-Trichlorophenol	0.360	0.374	-3.9	101	0.00
44		2,4,5-Trichlorophenol	0.380	0.395	-3.9	102	0.00

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Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Feb 03 15:16:18 2020
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)
45	S	2-Fluorobiphenyl	1.377	1.398	-1.5	97	0.00
46		1,1'-Biphenyl	1.559	1.562	-0.2	98	0.00
47		2-Chloronaphthalene	1.228	1.224	0.3	97	0.00
48		2-Nitroaniline	0.279	0.300	-7.5	104	0.00
49		Acenaphthylene	1.852	1.855	-0.2	98	0.00
50		Dimethylphthalate	1.476	1.489	-0.9	99	0.00
51		2,6-Dinitrotoluene	0.283	0.300	-6.0	103	0.00
52	C	Acenaphthene	1.186	1.179	0.6	98	0.00
53		3-Nitroaniline	0.333	0.351	-5.4	103	0.00
54	P	2,4-Dinitrophenol	0.078	0.086	-10.3	123	0.00
55		Dibenzofuran	1.755	1.756	-0.1	99	0.00
56	P	4-Nitrophenol	0.241	0.254	-5.4	103	0.00
57		2,4-Dinitrotoluene	0.369	0.396	-7.3	105	0.00
58		Fluorene	1.409	1.409	0.0	96	0.00
59		2,3,4,6-Tetrachlorophenol	0.330	0.350	-6.1	102	0.00
60		Diethylphthalate	1.477	1.481	-0.3	99	0.00
61		4-Chlorophenyl-phenylether	0.716	0.725	-1.3	99	0.00
62		4-Nitroaniline	0.351	0.376	-7.1	103	0.00
63		Azobenzene	1.352	1.332	1.5	97	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	100	0.00
65		4,6-Dinitro-2-methylphenol	0.057	0.061	-7.0	116	0.00
66	c	n-Nitrosodiphenylamine	0.596	0.591	0.8	98	0.00
67		4-Bromophenyl-phenylether	0.217	0.214	1.4	100	0.00
68		Hexachlorobenzene	0.247	0.241	2.4	99	0.00
69		Atrazine	0.179	0.184	-2.8	99	0.00
70	C	Pentachlorophenol	0.115	0.120	-4.3	104	0.00
71		Phenanthrene	1.077	1.064	1.2	98	0.00
72		Anthracene	1.042	1.036	0.6	98	0.00
73		Carbazole	0.993	0.989	0.4	99	0.00
74		Di-n-butylphthalate	1.160	1.190	-2.6	101	0.00
75	C	Fluoranthene	1.210	1.240	-2.5	102	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	103	0.00
77		Benzidine	0.470	0.543	-15.5	114	0.00
78		Pyrene	1.376	1.387	-0.8	102	0.00
79	S	Terphenyl-d14	0.980	0.980	0.0	97	0.00
80		Butylbenzylphthalate	0.540	0.559	-3.5	107	0.00
81		Benzo(a)anthracene	1.331	1.356	-1.9	105	0.01
82		3,3'-Dichlorobenzidine	0.449	0.484	-7.8	109	0.00
83		Chrysene	1.255	1.262	-0.6	102	0.01
84		Bis(2-ethylhexyl)phthalate	0.848	0.901	-6.3	109	0.00
85	c	Di-n-octyl phthalate	1.334	1.444	-8.2	111	0.00
86		Indeno(1,2,3-cd)pyrene	1.592	1.653	-3.8	108	0.02
87	I	Perylene-d12	1.000	1.000	0.0	109	0.01

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88	Benzo(b)fluoranthene	1.212	1.209	0.2	106	0.01
89	Benzo(k)fluoranthene	1.142	1.110	2.8	105	0.02
90 C	Benzo(a)pyrene	1.106	1.099	0.6	109	0.02
91	Dibenzo(a,h)anthracene	1.130	1.140	-0.9	109	0.02
92	Benzo(g,h,i)perylene	1.115	1.116	-0.1	110	0.02

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