Evaluate Continuing Calibration Report

Data Path : Z:\SVOASRV\HPCHEM1\BNA P\DATA\BP011520\
Data File : BP001606.D
Acq On : 15 Jan 2020 13:19
Operator : JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_P
LabSampleId : SSTDCCC040

Quant Time: Jan 16 06:21:38 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA P\METHODS\8270-BP010820.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jan 14 21:52:04 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	150#	0.00
2		1,4-Dioxane	0.326	0.393	-20.6	192#	0.00
3		Pyridine	1.177	1.216	-3.3	165#	0.00
4		n-Nitrosodimethylamine	0.834	0.979	-17.4	189#	0.00
5	S	2-Fluorophenol	1.021	0.999	2.2	157#	0.00
6		Aniline	2.040	1.915	6.1	148	0.00
7	S	Phenol-d6	1.632	1.510	7.5	150#	0.00
8		2-Chlorophenol	1.192	1.142	4.2	157#	0.00
9		Benzaldehyde	0.975	0.886	9.1	160#	0.00
10	C	Phenol	1.688	1.556	7.8	148	0.00
11		bis(2-Chloroethyl)ether	1.324	1.231	7.0	151#	0.00
12		1,3-Dichlorobenzene	1.498	1.521	-1.5	161#	0.00
13	C	1,4-Dichlorobenzene	1.549	1.550	-0.1	161#	0.00
14		1,2-Dichlorobenzene	1.497	1.485	0.8	161#	0.00
15		Benzyl Alcohol	1.537	1.302	15.3	139	0.00
16		2,2'-oxybis(1-Chloropropane	2.459	2.498	-1.6	160#	0.00
17		2-Methylphenol	1.172	1.071	8.6	149	0.00
18		Hexachloroethane	0.605	0.621	-2.6	165#	0.00
19	P	n-Nitroso-di-n-propylamine	1.244	1.130	9.2	146	0.00
20		3+4-Methylphenols	1.639	1.448	11.7	143	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	134	0.00
22		Acetophenone	0.561	0.561	0.0	143	0.00
23	S	Nitrobenzene-d5	0.458	0.479	-4.6	151#	0.00
24		Nitrobenzene	0.463	0.472	-1.9	148	0.00
25		Isophorone	0.821	0.790	3.8	139	0.00
26	C	2-Nitrophenol	0.175	0.173	1.1	144	0.00
27		2,4-Dimethylphenol	0.261	0.255	2.3	141	0.00
28		bis(2-Chloroethoxy)methane	0.481	0.473	1.7	142	0.00
29	C	2,4-Dichlorophenol	0.358	0.343	4.2	139	0.00
30		1,2,4-Trichlorobenzene	0.432	0.444	-2.8	148	0.00
31		Naphthalene	1.056	1.058	-0.2	145	0.00
32		Benzoic acid	0.222	0.168	24.3	109	0.02
33		4-Chloroaniline	0.486	0.429	11.7	130	0.00
34	C	Hexachlorobutadiene	0.306	0.325	-6.2	151#	0.00
35		Caprolactam	0.130	0.109	16.2	125	0.00
36	C	4-Chloro-3-methylphenol	0.429	0.372	13.3	127	0.00
37		2-Methylnaphthalene	0.841	0.781	7.1	135	0.00
38		1-Methylnaphthalene	0.810	0.756	6.7	137	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	118	0.00
40		1,2,4,5-Tetrachlorobenzene	0.685	0.759	-10.8	139	0.00
41	P	Hexachlorocyclopentadiene	0.348	0.349	-0.3	122	0.00
42	S	2,4,6-Tribromophenol	0.304	0.302	0.7	127	0.00
43	C	2,4,6-Trichlorophenol	0.442	0.453	-2.5	130	0.00
44		2,4,5-Trichlorophenol	0.470	0.478	-1.7	129	0.00

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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	1.360	1.484	-9.1	137	0.00
46		1,1'-Biphenyl	1.433	1.523	-6.3	134	0.00
47		2-Chloronaphthalene	1.191	1.256	-5.5	134	0.00
48		2-Nitroaniline	0.468	0.476	-1.7	129	0.00
49		Acenaphthylene	1.837	1.869	-1.7	129	0.00
50		Dimethylphthalate	1.647	1.582	3.9	124	0.00
51		2,6-Dinitrotoluene	0.349	0.336	3.7	124	0.00
52	C	Acenaphthene	1.138	1.119	1.7	127	0.00
53		3-Nitroaniline	0.369	0.344	6.8	122	0.00
54	P	2,4-Dinitrophenol	0.196	0.224	-14.3	149	0.00
55		Dibenzofuran	1.924	1.894	1.6	125	0.00
56	P	4-Nitrophenol	0.295	0.352	-19.3	155#	0.00
57		2,4-Dinitrotoluene	0.511	0.514	-0.6	129	0.00
58		Fluorene	1.557	1.533	1.5	126	0.00
59		2,3,4,6-Tetrachlorophenol	0.485	0.472	2.7	126	0.00
60		Diethylphthalate	1.694	1.710	-0.9	130	0.00
61		4-Chlorophenyl-phenylether	0.891	0.900	-1.0	126	0.00
62		4-Nitroaniline	0.421	0.381	9.5	119	0.00
63		Azobenzene	1.693	1.749	-3.3	133	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	118	0.00
65		4,6-Dinitro-2-methylphenol	0.108	0.111	-2.8	129	0.00
66	c	n-Nitrosodiphenylamine	0.530	0.529	0.2	126	0.00
67		4-Bromophenyl-phenylether	0.226	0.223	1.3	124	0.00
68		Hexachlorobenzene	0.262	0.265	-1.1	128	0.00
69		Atrazine	0.195	0.165	15.4	102	0.00
70	C	Pentachlorophenol	0.145	0.166	-14.5	147	0.00
71		Phenanthrene	1.088	1.087	0.1	128	0.00
72		Anthracene	1.059	1.088	-2.7	131	0.00
73		Carbazole	0.994	0.964	3.0	124	0.00
74		Di-n-butylphthalate	1.176	1.202	-2.2	130	0.00
75	C	Fluoranthene	1.431	1.408	1.6	128	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	119	0.00
77		Benzidine	0.457	0.482	-5.5	143	0.00
78		Pyrene	1.469	1.481	-0.8	128	0.00
79	S	Terphenyl-d14	1.093	1.105	-1.1	127	0.00
80		Butylbenzylphthalate	0.558	0.563	-0.9	127	0.00
81		Benzo(a)anthracene	1.393	1.362	2.2	125	0.00
82		3,3'-Dichlorobenzidine	0.522	0.482	7.7	119	0.00
83		Chrysene	1.357	1.354	0.2	129	0.00
84		Bis(2-ethylhexyl)phthalate	0.765	0.806	-5.4	133	0.00
85	c	Di-n-octyl phthalate	1.221	1.282	-5.0	134	0.00
86		Indeno(1,2,3-cd)pyrene	1.589	1.342	15.5	112	0.00
87	I	Perylene-d12	1.000	1.000	0.0	112	0.00

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	Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
88	Benzo(b)fluoranthene	1.260	1.297	-2.9	125	0.00
89	Benzo(k)fluoranthene	1.229	1.298	-5.6	128	0.00
90 C	Benzo(a)pyrene	1.197	1.203	-0.5	122	0.00
91	Dibenzo(a,h)anthracene	1.235	1.122	9.1	111	0.00
92	Benzo(g,h,i)perylene	1.227	1.100	10.4	111	0.00

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