Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP040122\
Data File : BP009652.D
Acq On : 01 Apr 2022 11:40
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Apr 01 22:42:32 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP031722.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Mar 31 04:36:51 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	105	0.01
2		1,4-Dioxane	0.454	0.489	-7.7	119	0.00
3		Pyridine	1.222	1.361	-11.4	120	0.00
4		n-Nitrosodimethylamine	0.450	0.503	-11.8	122	0.00
5	S	2-Fluorophenol	1.146	1.257	-9.7	119	0.00
6		Aniline	1.766	1.868	-5.8	115	0.00
7	S	Phenol-d6	1.549	1.659	-7.1	117	0.01
8		2-Chlorophenol	1.306	1.355	-3.8	113	0.00
9		Benzaldehyde	0.821	0.825	-0.5	125	0.00
10	C	Phenol	1.554	1.667	-7.3	117	0.01
11		bis(2-Chloroethyl)ether	1.230	1.310	-6.5	117	0.00
12		1,3-Dichlorobenzene	1.474	1.508	-2.3	112	0.01
13	C	1,4-Dichlorobenzene	1.507	1.537	-2.0	112	0.01
14		1,2-Dichlorobenzene	1.457	1.474	-1.2	111	0.01
15		Benzyl Alcohol	1.235	1.255	-1.6	111	0.00
16		2,2'-oxybis(1-Chloropropane	1.333	1.511	-13.4	124	0.01
17		2-Methylphenol	1.071	1.104	-3.1	113	0.00
18		Hexachloroethane	0.528	0.545	-3.2	113	0.01
19	P	n-Nitroso-di-n-propylamine	0.980	0.987	-0.7	111	0.00
20		3+4-Methylphenols	1.473	1.507	-2.3	111	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	103	0.01
22		Acetophenone	0.495	0.517	-4.4	111	0.00
23	S	Nitrobenzene-d5	0.376	0.405	-7.7	113	0.00
24		Nitrobenzene	0.356	0.383	-7.6	114	0.00
25		Isophorone	0.642	0.674	-5.0	112	0.01
26	C	2-Nitrophenol	0.185	0.195	-5.4	110	0.00
27		2,4-Dimethylphenol	0.261	0.279	-6.9	112	0.01
28		bis(2-Chloroethoxy)methane	0.418	0.446	-6.7	113	0.01
29	C	2,4-Dichlorophenol	0.319	0.333	-4.4	109	0.01
30		1,2,4-Trichlorobenzene	0.370	0.375	-1.4	107	0.01
31		Naphthalene	1.030	1.054	-2.3	109	0.01
32		Benzoic acid	0.204	0.223	-9.3	112	0.01
33		4-Chloroaniline	0.420	0.431	-2.6	108	0.01
34	C	Hexachlorobutadiene	0.250	0.240	4.0	101	0.01
35		Caprolactam	0.108	0.108	0.0	110	0.01
36	C	4-Chloro-3-methylphenol	0.343	0.338	1.5	106	0.00
37		2-Methylnaphthalene	0.723	0.726	-0.4	108	0.01
38		1-Methylnaphthalene	0.704	0.705	-0.1	107	0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	99	0.01
40		1,2,4,5-Tetrachlorobenzene	0.664	0.679	-2.3	103	0.01
41	P	Hexachlorocyclopentadiene	0.234	0.262	-12.0	107	0.02
42	S	2,4,6-Tribromophenol	0.330	0.292	11.5	90	0.01
43	C	2,4,6-Trichlorophenol	0.440	0.458	-4.1	104	0.01
44		2,4,5-Trichlorophenol	0.478	0.478	0.0	101	0.00
45	S	2-Fluorobiphenyl	1.443	1.473	-2.1	103	0.01
46		1,1'-Biphenyl	1.492	1.559	-4.5	106	0.01
47		2-Chloronaphthalene	1.175	1.228	-4.5	106	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.320	0.357	-11.6	112	0.00
49		Acenaphthylene	1.813	1.886	-4.0	106	0.01
50		Dimethylphthalate	1.515	1.519	-0.3	103	0.01
51		2,6-Dinitrotoluene	0.318	0.323	-1.6	102	0.00
52	C	Acenaphthene	1.083	1.117	-3.1	105	0.01
53		3-Nitroaniline	0.335	0.354	-5.7	106	0.00
54	P	2,4-Dinitrophenol	0.174	0.167	4.0	95	0.00
55		Dibenzofuran	1.820	1.854	-1.9	104	0.01
56	P	4-Nitrophenol	0.250	0.252	-0.8	99	0.00
57		2,4-Dinitrotoluene	0.437	0.437	0.0	100	0.01
58		Fluorene	1.431	1.441	-0.7	104	0.02
59		2,3,4,6-Tetrachlorophenol	0.433	0.414	4.4	97	0.01
60		Diethylphthalate	1.517	1.503	0.9	102	0.01
61		4-Chlorophenyl-phenylether	0.788	0.773	1.9	101	0.01
62		4-Nitroaniline	0.352	0.372	-5.7	106	0.01
63		Azobenzene	1.319	1.418	-7.5	109	0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	95	0.01
65		4,6-Dinitro-2-methylphenol	0.129	0.136	-5.4	97	0.01
66	c	n-Nitrosodiphenylamine	0.592	0.624	-5.4	103	0.01
67		4-Bromophenyl-phenylether	0.245	0.241	1.6	97	0.01
68		Hexachlorobenzene	0.282	0.268	5.0	93	0.01
69		Atrazine	0.208	0.211	-1.4	96	0.01
70	C	Pentachlorophenol	0.151	0.140	7.3	86	0.01
71		Phenanthrene	1.071	1.105	-3.2	102	0.02
72		Anthracene	1.057	1.106	-4.6	103	0.01
73		Carbazole	1.034	1.086	-5.0	103	0.01
74		Di-n-butylphthalate	1.211	1.255	-3.6	101	0.01
75	C	Fluoranthene	1.296	1.330	-2.6	102	0.01
76	I	Chrysene-d12	1.000	1.000	0.0	100	0.02
77		Benzidine	0.490	0.525	-7.1	113	0.01
78		Pyrene	1.318	1.332	-1.1	102	0.01
79	S	Terphenyl-d14	1.114	1.067	4.2	96	0.01
80		Butylbenzylphthalate	0.560	0.576	-2.9	104	0.01
81		Benzo(a)anthracene	1.314	1.335	-1.6	105	0.01
82		3,3'-Dichlorobenzidine	0.508	0.516	-1.6	105	0.01
83		Chrysene	1.244	1.278	-2.7	106	0.01
84		Bis(2-ethylhexyl)phthalate	0.781	0.812	-4.0	106	0.01
85	c	Di-n-octyl phthalate	1.346	1.435	-6.6	110	0.02
86	I	Perylene-d12	1.000	1.000	0.0	101	0.02
87		Indeno(1,2,3-cd)pyrene	1.517	1.530	-0.9	105	0.04
88		Benzo(b)fluoranthene	1.194	1.208	-1.2	106	0.02
89		Benzo(k)fluoranthene	1.169	1.238	-5.9	110	0.02
90	C	Benzo(a)pyrene	1.022	1.054	-3.1	108	0.02
91		Dibenzo(a,h)anthracene	1.265	1.277	-0.9	105	0.04
92		Benzo(g,h,i)perylene	1.244	1.257	-1.0	105	0.04

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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