Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080922\
Data File : BM036168.D
Acq On : 09 Aug 2022 10:28
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Aug 10 03:29:36 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM080122.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 08 23:36:56 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	88	0.00
2		1,4-Dioxane	0.497	0.475	4.4	95	0.00
3		Pyridine	1.333	1.342	-0.7	102	0.00
4		n-Nitrosodimethylamine	0.548	0.529	3.5	96	0.00
5	S	2-Fluorophenol	1.234	1.208	2.1	95	0.01
6		Aniline	1.738	1.900	-9.3	107	0.00
7	S	Phenol-d6	1.570	1.695	-8.0	106	0.02
8		2-Chlorophenol	1.315	1.372	-4.3	100	0.01
9		Benzaldehyde	0.831	0.830	0.1	110	0.01
10	C	Phenol	1.580	1.694	-7.2	105	0.02
11		bis(2-Chloroethyl)ether	1.302	1.319	-1.3	97	0.01
12		1,3-Dichlorobenzene	1.458	1.446	0.8	95	0.01
13	C	1,4-Dichlorobenzene	1.486	1.484	0.1	95	0.00
14		1,2-Dichlorobenzene	1.434	1.452	-1.3	96	0.00
15		Benzyl Alcohol	1.053	1.224	-16.2	113	0.01
16		2,2'-oxybis(1-Chloropropane	1.731	1.762	-1.8	96	0.01
17		2-Methylphenol	1.046	1.169	-11.8	110	0.02
18		Hexachloroethane	0.535	0.537	-0.4	93	0.00
19	P	n-Nitroso-di-n-propylamine	0.953	1.088	-14.2	109	0.00
20		3+4-Methylphenols	1.421	1.640	-15.4	112	0.01
21	I	Naphthalene-d8	1.000	1.000	0.0	99	0.00
22		Acetophenone	0.473	0.460	2.7	109	0.00
23	S	Nitrobenzene-d5	0.357	0.343	3.9	107	0.01
24		Nitrobenzene	0.336	0.323	3.9	108	0.01
25		Isophorone	0.581	0.604	-4.0	115	0.01
26	C	2-Nitrophenol	0.158	0.171	-8.2	118	0.01
27		2,4-Dimethylphenol	0.245	0.252	-2.9	114	0.02
28		bis(2-Chloroethoxy)methane	0.414	0.401	3.1	106	0.01
29	C	2,4-Dichlorophenol	0.276	0.286	-3.6	115	0.01
30		1,2,4-Trichlorobenzene	0.313	0.295	5.8	104	0.01
31		Naphthalene	0.994	0.945	4.9	107	0.01
32		Benzoic acid	0.195	0.251	-28.7#	144	0.05
33		4-Chloroaniline	0.401	0.417	-4.0	118	0.01
34	C	Hexachlorobutadiene	0.184	0.171	7.1	100	0.01
35		Caprolactam	0.085	0.107	-25.9#	149	0.03
36	C	4-Chloro-3-methylphenol	0.290	0.323	-11.4	127	0.02
37		2-Methylnaphthalene	0.662	0.667	-0.8	113	0.01
38		1-Methylnaphthalene	0.648	0.653	-0.8	114	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	113	0.01
40		1,2,4,5-Tetrachlorobenzene	0.542	0.487	10.1	112	0.00
41	P	Hexachlorocyclopentadiene	0.287	0.310	-8.0	128	0.01
42	S	2,4,6-Tribromophenol	0.235	0.245	-4.3	134	0.01
43	C	2,4,6-Trichlorophenol	0.369	0.366	0.8	124	0.01
44		2,4,5-Trichlorophenol	0.389	0.393	-1.0	127	0.01
45	S	2-Fluorobiphenyl	1.355	1.218	10.1	113	0.01
46		1,1'-Biphenyl	1.458	1.330	8.8	117	0.01
47		2-Chloronaphthalene	1.116	1.018	8.8	117	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.304	0.323	-6.3	140	0.01
49		Acenaphthylene	1.724	1.656	3.9	124	0.01
50		Dimethylphthalate	1.360	1.336	1.8	128	0.01
51		2,6-Dinitrotoluene	0.273	0.285	-4.4	133	0.01
52	C	Acenaphthene	1.128	1.104	2.1	127	0.01
53		3-Nitroaniline	0.318	0.349	-9.7	144	0.02
54	P	2,4-Dinitrophenol	0.144	0.183	-27.1#	165#	0.01
55		Dibenzofuran	1.689	1.613	4.5	125	0.01
56	P	4-Nitrophenol	0.255	0.303	-18.8	158#	0.02
57		2,4-Dinitrotoluene	0.358	0.404	-12.8	145	0.01
58		Fluorene	1.332	1.286	3.5	127	0.01
59		2,3,4,6-Tetrachlorophenol	0.330	0.339	-2.7	133	0.01
60		Diethylphthalate	1.398	1.389	0.6	128	0.01
61		4-Chlorophenyl-phenylether	0.664	0.631	5.0	122	0.01
62		4-Nitroaniline	0.328	0.379	-15.5	153#	0.02
63		Azobenzene	1.329	1.273	4.2	121	0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	125	0.01
65		4,6-Dinitro-2-methylphenol	0.102	0.115	-12.7	154#	0.02
66	c	n-Nitrosodiphenylamine	0.565	0.525	7.1	131	0.01
67		4-Bromophenyl-phenylether	0.200	0.184	8.0	128	0.01
68		Hexachlorobenzene	0.230	0.215	6.5	131	0.01
69		Atrazine	0.158	0.165	-4.4	137	0.01
70	C	Pentachlorophenol	0.134	0.136	-1.5	141	0.01
71		Phenanthrene	1.019	0.954	6.4	135	0.01
72		Anthracene	0.986	0.941	4.6	136	0.01
73		Carbazole	1.000	0.985	1.5	143	0.01
74		Di-n-butylphthalate	1.193	1.157	3.0	133	0.01
75	C	Fluoranthene	1.141	1.114	2.4	142	0.01
76	I	Chrysene-d12	1.000	1.000	0.0	131	0.02
77		Benzidine	0.300	0.475	-58.3#	217#	0.00
78		Pyrene	1.240	1.169	5.7	144	0.01
79	S	Terphenyl-d14	1.014	0.900	11.2	130	0.01
80		Butylbenzylphthalate	0.535	0.554	-3.6	149	0.01
81		Benzo(a)anthracene	1.232	1.173	4.8	144	0.01
82		3,3'-Dichlorobenzidine	0.299	0.317	-6.0	156#	0.01
83		Chrysene	1.171	1.117	4.6	144	0.01
84		Bis(2-ethylhexyl)phthalate	0.808	0.778	3.7	135	0.00
85	c	Di-n-octyl phthalate	1.318	1.347	-2.2	142	0.01
86	I	Perylene-d12	1.000	1.000	0.0	143	0.02
87		Indeno(1,2,3-cd)pyrene	1.405	1.372	2.3	160#	0.04
88		Benzo(b)fluoranthene	1.172	1.071	8.6	149	0.02
89		Benzo(k)fluoranthene	1.184	1.081	8.7	149	0.02
90	C	Benzo(a)pyrene	0.983	0.932	5.2	155#	0.02
91		Dibenzo(a,h)anthracene	1.176	1.141	3.0	159#	0.04
92		Benzo(g,h,i)perylene	1.139	1.129	0.9	164#	0.04

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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