Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG040924\
Data File : BG060837.D
Acq On : 9 Apr 2024 9:31
Operator : MA/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Apr 09 10:05:25 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG032924.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Sat Mar 30 03:12:26 2024
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	110	-0.01
2		1,4-Dioxane	0.492	0.477	3.0	110	0.00
3		Pyridine	1.485	1.365	8.1	93	0.00
4		n-Nitrosodimethylamine	0.887	0.875	1.4	103	0.00
5	S	2-Fluorophenol	1.201	1.130	5.9	103	0.00
6		Aniline	2.250	1.956	13.1	94	-0.01
7	S	Phenol-d6	1.782	1.651	7.4	100	0.00
8		2-Chlorophenol	1.361	1.265	7.1	101	-0.01
9		Benzaldehyde	0.983	0.858	12.7	99	-0.01
10	C	Phenol	1.819	1.666	8.4	97	0.00
11		bis(2-Chloroethyl)ether	1.468	1.258	14.3	94	-0.01
12		1,3-Dichlorobenzene	1.415	1.347	4.8	106	-0.01
13	C	1,4-Dichlorobenzene	1.444	1.355	6.2	104	-0.01
14		1,2-Dichlorobenzene	1.415	1.320	6.7	104	-0.01
15		Benzyl Alcohol	1.444	1.302	9.8	96	0.00
16		2,2'-oxybis(1-Chloropropane	3.314	2.851	14.0	95	-0.01
17		2-Methylphenol	1.373	1.218	11.3	96	0.00
18		Hexachloroethane	0.513	0.470	8.4	104	-0.01
19	P	n-Nitroso-di-n-propylamine	1.359	1.124	17.3	89	-0.01
20		3+4-Methylphenols	1.880	1.623	13.7	93	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	94	-0.01
22		Acetophenone	0.551	0.522	5.3	91	-0.01
23	S	Nitrobenzene-d5	0.350	0.400	-14.3	110	-0.01
24		Nitrobenzene	0.352	0.395	-12.2	102	-0.01
25		Isophorone	0.807	0.735	8.9	86	-0.01
26	C	2-Nitrophenol	0.120	0.160	-33.3#	121	-0.01
27		2,4-Dimethylphenol	0.324	0.300	7.4	88	-0.01
28		bis(2-Chloroethoxy)methane	0.477	0.436	8.6	87	-0.02
29	C	2,4-Dichlorophenol	0.298	0.318	-6.7	98	-0.01
30		1,2,4-Trichlorobenzene	0.337	0.348	-3.3	99	-0.02
31		Naphthalene	1.082	1.044	3.5	91	-0.02
32		Benzoic acid	0.191	0.218	-14.1	103	0.00
33		4-Chloroaniline	0.503	0.478	5.0	88	-0.01
34	C	Hexachlorobutadiene	0.225	0.245	-8.9	105	-0.02
35		Caprolactam	0.118	0.115	2.5	89	-0.01
36	C	4-Chloro-3-methylphenol	0.380	0.381	-0.3	91	0.00
37		2-Methylnaphthalene	0.795	0.767	3.5	92	-0.01
38		1-Methylnaphthalene	0.751	0.728	3.1	94	-0.01
39	I	Acenaphthene-d10	1.000	1.000	0.0	95	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.575	0.619	-7.7	106	-0.01
41	P	Hexachlorocyclopentadiene	0.292	0.321	-9.9	106	-0.01
42	S	2,4,6-Tribromophenol	0.227	0.294	-29.5#	123	0.00
43	C	2,4,6-Trichlorophenol	0.370	0.415	-12.2	106	0.00
44		2,4,5-Trichlorophenol	0.442	0.501	-13.3	107	0.00
45	S	2-Fluorobiphenyl	1.363	1.379	-1.2	99	-0.02
46		1,1'-Biphenyl	1.413	1.372	2.9	96	-0.01
47		2-Chloronaphthalene	1.073	1.047	2.4	95	-0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.307	0.405	-31.9#	119	0.00
49		Acenaphthylene	1.803	1.753	2.8	95	-0.01
50		Dimethylphthalate	1.581	1.523	3.7	92	-0.02
51		2,6-Dinitrotoluene	0.247	0.307	-24.3	110	-0.01
52	C	Acenaphthene	1.132	1.123	0.8	96	-0.01
53		3-Nitroaniline	0.277	0.327	-18.1	104	0.00
54	P	2,4-Dinitrophenol	0.110	0.160	-45.5#	141	0.00
55		Dibenzofuran	1.786	1.712	4.1	94	-0.01
56	P	4-Nitrophenol	0.231	0.271	-17.3	109	0.00
57		2,4-Dinitrotoluene	0.320	0.434	-35.6#	119	0.00
58		Fluorene	1.433	1.408	1.7	95	-0.01
59		2,3,4,6-Tetrachlorophenol	0.385	0.442	-14.8	105	-0.01
60		Diethylphthalate	1.590	1.488	6.4	91	-0.02
61		4-Chlorophenyl-phenylether	0.758	0.753	0.7	97	-0.02
62		4-Nitroaniline	0.289	0.349	-20.8	110	0.00
63		Azobenzene	1.550	1.401	9.6	87	-0.01
64	I	Phenanthrene-d10	1.000	1.000	0.0	98	-0.01
65		4,6-Dinitro-2-methylphenol	0.073	0.103	-41.1#	139	0.00
66	c	n-Nitrosodiphenylamine	0.572	0.546	4.5	94	-0.01
67		4-Bromophenyl-phenylether	0.203	0.216	-6.4	105	-0.02
68		Hexachlorobenzene	0.229	0.236	-3.1	102	-0.01
69		Atrazine	0.200	0.202	-1.0	99	-0.02
70	C	Pentachlorophenol	0.150	0.173	-15.3	111	0.00
71		Phenanthrene	1.040	1.009	3.0	98	-0.01
72		Anthracene	1.056	1.039	1.6	98	-0.02
73		Carbazole	0.931	0.907	2.6	98	0.00
74		Di-n-butylphthalate	1.161	1.082	6.8	92	-0.02
75	C	Fluoranthene	1.262	1.272	-0.8	102	-0.01
76	I	Chrysene-d12	1.000	1.000	0.0	107	-0.02
77		Benzidine	0.527	0.509	3.4	92	-0.01
78		Pyrene	1.399	1.317	5.9	103	-0.01
79	S	Terphenyl-d14	1.136	1.078	5.1	105	-0.02
80		Butylbenzylphthalate	0.510	0.484	5.1	100	-0.02
81		Benzo(a)anthracene	1.410	1.363	3.3	104	-0.02
82		3,3'-Dichlorobenzidine	0.476	0.482	-1.3	105	-0.02
83		Chrysene	1.359	1.314	3.3	105	-0.02
84		Bis(2-ethylhexyl)phthalate	0.813	0.722	11.2	93	-0.03
85	c	Di-n-octyl phthalate	1.325	1.241	6.3	96	-0.05
86	I	Perylene-d12	1.000	1.000	0.0	108	-0.03
87		Indeno(1,2,3-cd)pyrene	1.408	1.357	3.6	105	-0.05
88		Benzo(b)fluoranthene	1.149	1.105	3.8	104	-0.03
89		Benzo(k)fluoranthene	1.176	1.111	5.5	101	-0.02
90	C	Benzo(a)pyrene	1.117	1.088	2.6	104	-0.02
91		Dibenzo(a,h)anthracene	1.187	1.146	3.5	105	-0.05
92		Benzo(g,h,i)perylene	1.155	1.092	5.5	103	-0.04

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

(#) = Out of Range SPCC's out = 0 CCC's out = 1