Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG122222\
Data File : BG056059.D
Acq On : 22 Dec 2022 11:00
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_G
LabSampleId :
SSTDCCC040

Quant Time: Dec 23 05:23:21 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG121322.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Dec 23 05:09:54 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	180#	0.00
2		1,4-Dioxane	0.248	0.256	-3.2	180#	0.00
3		Pyridine	0.817	0.879	-7.6	183#	0.00
4		n-Nitrosodimethylamine	0.734	0.689	6.1	168#	0.00
5	S	2-Fluorophenol	0.853	0.891	-4.5	189#	0.00
6		Aniline	1.313	1.397	-6.4	186#	0.00
7	S	Phenol-d6	1.097	1.197	-9.1	198#	0.00
8		2-Chlorophenol	1.100	1.175	-6.8	192#	0.00
9		Benzaldehyde	0.689	0.669	2.9	176#	0.00
10	C	Phenol	1.200	1.291	-7.6	193#	0.00
11		bis(2-Chloroethyl)ether	0.782	0.847	-8.3	197#	0.00
12		1,3-Dichlorobenzene	1.423	1.410	0.9	179#	0.00
13	C	1,4-Dichlorobenzene	1.455	1.476	-1.4	189#	0.00
14		1,2-Dichlorobenzene	1.390	1.413	-1.7	182#	0.00
15		Benzyl Alcohol	1.140	1.154	-1.2	184#	0.00
16		2,2'-oxybis(1-Chloropropane	1.018	1.245	-22.3	227#	0.00
17		2-Methylphenol	0.944	0.995	-5.4	187#	0.00
18		Hexachloroethane	0.442	0.469	-6.1	186#	0.00
19	P	n-Nitroso-di-n-propylamine	0.786	0.885	-12.6	205#	0.00
20		3+4-Methylphenols	1.348	1.440	-6.8	192#	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	194#	0.00
22		Acetophenone	0.496	0.483	2.6	185#	0.00
23	S	Nitrobenzene-d5	0.286	0.339	-18.5	226#	0.00
24		Nitrobenzene	0.288	0.336	-16.7	216#	0.00
25		Isophorone	0.600	0.623	-3.8	196#	0.00
26	C	2-Nitrophenol	0.171	0.198	-15.8	219#	0.00
27		2,4-Dimethylphenol	0.286	0.271	5.2	179#	0.00
28		bis(2-Chloroethoxy)methane	0.265	0.287	-8.3	215#	0.00
29	C	2,4-Dichlorophenol	0.407	0.420	-3.2	199#	0.00
30		1,2,4-Trichlorobenzene	0.528	0.519	1.7	192#	0.00
31		Naphthalene	1.055	1.040	1.4	193#	0.00
32		Benzoic acid	0.221	0.260	-17.6	208#	0.00
33		4-Chloroaniline	0.452	0.454	-0.4	188#	0.00
34	C	Hexachlorobutadiene	0.497	0.446	10.3	183#	0.00
35		Caprolactam	0.113	0.120	-6.2	203#	0.00
36	C	4-Chloro-3-methylphenol	0.381	0.410	-7.6	198#	0.00
37		2-Methylnaphthalene	0.932	0.938	-0.6	194#	0.00
38		1-Methylnaphthalene	0.908	0.903	0.6	193#	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	192#	0.00
40		1,2,4,5-Tetrachlorobenzene	0.844	0.823	2.5	184#	0.00
41	P	Hexachlorocyclopentadiene	0.479	0.479	0.0	190#	0.00
42	S	2,4,6-Tribromophenol	0.482	0.520	-7.9	192#	0.00
43	C	2,4,6-Trichlorophenol	0.500	0.551	-10.2	204#	0.00
44		2,4,5-Trichlorophenol	0.563	0.611	-8.5	200#	0.00
45	S	2-Fluorobiphenyl	1.426	1.424	0.1	190#	0.00
46		1,1'-Biphenyl	1.356	1.416	-4.4	198#	0.00
47		2-Chloronaphthalene	1.102	1.132	-2.7	193#	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.184	0.300	-63.0#	271#	0.00
49		Acenaphthylene	1.758	1.806	-2.7	195#	0.00
50		Dimethylphthalate	1.729	1.763	-2.0	188#	0.00
51		2,6-Dinitrotoluene	0.268	0.358	-33.6#	223#	0.00
52	C	Acenaphthene	1.180	1.267	-7.4	201#	0.00
53		3-Nitroaniline	0.272	0.326	-19.9	216#	0.00
54	P	2,4-Dinitrophenol	0.199	0.244	-22.6	224#	0.00
55		Dibenzofuran	1.985	2.046	-3.1	190#	0.00
56	P	4-Nitrophenol	0.218	0.263	-20.6	211#	0.00
57		2,4-Dinitrotoluene	0.404	0.530	-31.2#	235#	0.00
58		Fluorene	1.623	1.669	-2.8	192#	0.00
59		2,3,4,6-Tetrachlorophenol	0.643	0.696	-8.2	190#	0.00
60		Diethylphthalate	1.690	1.707	-1.0	186#	0.00
61		4-Chlorophenyl-phenylether	1.105	1.121	-1.4	187#	0.00
62		4-Nitroaniline	0.296	0.354	-19.6	210#	0.00
63		Azobenzene	1.029	1.078	-4.8	196#	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	187#	0.00
65		4,6-Dinitro-2-methylphenol	0.096	0.127	-32.3#	238#	0.00
66	c	n-Nitrosodiphenylamine	0.524	0.528	-0.8	189#	0.00
67		4-Bromophenyl-phenylether	0.293	0.296	-1.0	191#	0.00
68		Hexachlorobenzene	0.361	0.348	3.6	181#	0.00
69		Atrazine	0.237	0.230	3.0	182#	0.00
70	C	Pentachlorophenol	0.206	0.217	-5.3	190#	0.00
71		Phenanthrene	1.048	1.066	-1.7	189#	0.00
72		Anthracene	1.037	1.033	0.4	186#	0.00
73		Carbazole	0.855	0.839	1.9	185#	0.00
74		Di-n-butylphthalate	0.995	0.959	3.6	180#	0.00
75	C	Fluoranthene	1.468	1.363	7.2	174#	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	161#	0.00
77		Benzidine	0.351	0.363	-3.4	163#	0.00
78		Pyrene	1.210	1.383	-14.3	173#	0.00
79	S	Terphenyl-d14	1.132	1.220	-7.8	164#	0.00
80		Butylbenzylphthalate	0.343	0.400	-16.6	174#	0.00
81		Benzo(a)anthracene	1.407	1.490	-5.9	161#	0.00
82		3,3'-Dichlorobenzidine	0.506	0.540	-6.7	160#	0.00
83		Chrysene	1.319	1.416	-7.4	162#	0.00
84		Bis(2-ethylhexyl)phthalate	0.509	0.552	-8.4	163#	0.00
85	c	Di-n-octyl phthalate	0.864	0.934	-8.1	162#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	155#	0.00
87		Indeno(1,2,3-cd)pyrene	1.608	1.598	0.6	155#	0.00
88		Benzo(b)fluoranthene	1.294	1.300	-0.5	156#	0.00
89		Benzo(k)fluoranthene	1.291	1.301	-0.8	157#	0.00
90	C	Benzo(a)pyrene	1.094	1.104	-0.9	157#	0.00
91		Dibenzo(a,h)anthracene	1.346	1.341	0.4	156#	0.00
92		Benzo(g,h,i)perylene	1.301	1.307	-0.5	158#	0.00

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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