Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_G\Data\BG081721\
Data File : BG049871.D
Acq On : 17 Aug 2021 10:11
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_G
LabSampleId : SSTDCCC040

Quant Time: Aug 17 11:57:22 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_G\Methods\8270-BG080321.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Aug 16 18:23:02 2021
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	0.00
2		1,4-Dioxane	0.488	0.438	10.2	91	0.00
3		Pyridine	1.336	1.097	17.9	80	0.00
4		n-Nitrosodimethylamine	0.719	0.618	14.0	85	-0.01
5	S	2-Fluorophenol	1.116	1.006	9.9	88	0.00
6		Aniline	1.790	1.722	3.8	94	0.00
7	S	Phenol-d6	1.689	1.592	5.7	91	0.00
8		2-Chlorophenol	1.216	1.177	3.2	98	0.00
9		Benzaldehyde	1.127	1.098	2.6	98	0.00
10	C	Phenol	1.636	1.514	7.5	89	0.00
11		bis(2-Chloroethyl)ether	1.105	1.072	3.0	96	0.00
12		1,3-Dichlorobenzene	1.393	1.281	8.0	93	0.00
13	C	1,4-Dichlorobenzene	1.410	1.297	8.0	92	0.00
14		1,2-Dichlorobenzene	1.336	1.253	6.2	92	0.00
15		Benzyl Alcohol	1.418	1.392	1.8	94	0.00
16		2,2'-oxybis(1-Chloropropane	1.255	1.168	6.9	91	-0.01
17		2-Methylphenol	1.078	1.083	-0.5	96	0.00
18		Hexachloroethane	0.542	0.507	6.5	90	0.00
19	P	n-Nitroso-di-n-propylamine	1.130	1.115	1.3	96	0.00
20		3+4-Methylphenols	1.513	1.495	1.2	96	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	112	0.00
22		Acetophenone	0.543	0.467	14.0	95	0.00
23	S	Nitrobenzene-d5	0.452	0.393	13.1	94	0.00
24		Nitrobenzene	0.477	0.402	15.7	92	0.00
25		Isophorone	0.807	0.708	12.3	96	0.00
26	C	2-Nitrophenol	0.182	0.167	8.2	98	-0.01
27		2,4-Dimethylphenol	0.269	0.250	7.1	101	0.00
28		bis(2-Chloroethoxy)methane	0.353	0.323	8.5	100	0.00
29	C	2,4-Dichlorophenol	0.330	0.292	11.5	97	0.00
30		1,2,4-Trichlorobenzene	0.363	0.331	8.8	102	0.00
31		Naphthalene	1.047	0.905	13.6	96	0.00
32		Benzoic acid	0.174	0.129	25.9#	77	0.00
33		4-Chloroaniline	0.414	0.382	7.7	100	0.00
34	C	Hexachlorobutadiene	0.264	0.233	11.7	98	-0.01
35		Caprolactam	0.102	0.088	13.7	96	0.00
36	C	4-Chloro-3-methylphenol	0.381	0.357	6.3	101	0.00
37		2-Methylnaphthalene	0.789	0.722	8.5	99	0.00
38		1-Methylnaphthalene	0.742	0.653	12.0	97	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	116	0.00
40		1,2,4,5-Tetrachlorobenzene	0.589	0.530	10.0	101	0.00
41	P	Hexachlorocyclopentadiene	0.243	0.248	-2.1	94	0.00
42	S	2,4,6-Tribromophenol	0.294	0.267	9.2	104	0.00
43	C	2,4,6-Trichlorophenol	0.397	0.368	7.3	101	0.00
44		2,4,5-Trichlorophenol	0.431	0.386	10.4	98	0.00
45	S	2-Fluorobiphenyl	1.374	1.201	12.6	98	0.00
46		1,1'-Biphenyl	1.424	1.272	10.7	101	0.00
47		2-Chloronaphthalene	1.138	0.996	12.5	98	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.418	0.367	12.2	97	0.00
49		Acenaphthylene	1.803	1.592	11.7	100	0.00
50		Dimethylphthalate	1.509	1.314	12.9	100	0.00
51		2,6-Dinitrotoluene	0.332	0.300	9.6	98	0.00
52	C	Acenaphthene	1.086	0.957	11.9	99	0.00
53		3-Nitroaniline	0.324	0.286	11.7	99	0.00
54	P	2,4-Dinitrophenol	0.189	0.172	9.0	102	0.00
55		Dibenzofuran	1.762	1.530	13.2	97	0.00
56	P	4-Nitrophenol	0.237	0.195	17.7	88	0.00
57		2,4-Dinitrotoluene	0.467	0.429	8.1	105	0.00
58		Fluorene	1.432	1.272	11.2	101	0.00
59		2,3,4,6-Tetrachlorophenol	0.400	0.336	16.0	96	0.00
60		Diethylphthalate	1.518	1.318	13.2	99	0.00
61		4-Chlorophenyl-phenylether	0.764	0.663	13.2	99	0.00
62		4-Nitroaniline	0.333	0.298	10.5	98	0.00
63		Azobenzene	1.606	1.350	15.9	96	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	116	0.00
65		4,6-Dinitro-2-methylphenol	0.129	0.123	4.7	104	0.00
66	c	n-Nitrosodiphenylamine	0.565	0.521	7.8	101	0.00
67		4-Bromophenyl-phenylether	0.236	0.216	8.5	101	0.00
68		Hexachlorobenzene	0.262	0.239	8.8	101	0.00
69		Atrazine	0.226	0.208	8.0	102	0.00
70	C	Pentachlorophenol	0.140	0.113	19.3	88	0.00
71		Phenanthrene	1.052	0.945	10.2	100	0.00
72		Anthracene	1.033	0.930	10.0	101	0.00
73		Carbazole	0.979	0.881	10.0	99	0.00
74		Di-n-butylphthalate	1.141	1.007	11.7	98	0.00
75	C	Fluoranthene	1.295	1.087	16.1	94	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	100	0.00
77		Benzidine	0.512	0.565	-10.4	94	0.00
78		Pyrene	1.363	1.313	3.7	93	0.00
79	S	Terphenyl-d14	1.101	1.060	3.7	93	0.00
80		Butylbenzylphthalate	0.519	0.472	9.1	87	0.00
81		Benzo(a)anthracene	1.302	1.167	10.4	87	0.00
82		3,3'-Dichlorobenzidine	0.380	0.343	9.7	85	0.00
83		Chrysene	1.245	1.124	9.7	89	0.00
84		Bis(2-ethylhexyl)phthalate	0.735	0.662	9.9	87	0.00
85	c	Di-n-octyl phthalate	1.242	1.099	11.5	85	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	97	0.00
87		Indeno(1,2,3-cd)pyrene	1.442	1.293	10.3	86	0.00
88		Benzo(b)fluoranthene	1.312	1.165	11.2	86	-0.01
89		Benzo(k)fluoranthene	1.224	1.084	11.4	84	0.00
90	C	Benzo(a)pyrene	1.187	1.060	10.7	86	-0.01
91		Dibenzo(a,h)anthracene	1.215	1.119	7.9	87	0.00
92		Benzo(g,h,i)perylene	1.191	1.077	9.6	87	0.00

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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