Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM072222\
Data File : BM035902.D
Acq On : 22 Jul 2022 12:10
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_M
LabSampleId :
SSTDCCC040

Quant Time: Jul 22 14:28:11 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM071822.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jul 22 14:25:12 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	112	0.00
2		1,4-Dioxane	0.533	0.534	-0.2	110	0.00
3		Pyridine	1.417	1.423	-0.4	109	0.00
4		n-Nitrosodimethylamine	0.592	0.591	0.2	108	0.00
5	S	2-Fluorophenol	1.290	1.312	-1.7	110	0.00
6		Aniline	1.782	1.799	-1.0	109	0.00
7	S	Phenol-d6	1.622	1.629	-0.4	110	0.00
8		2-Chlorophenol	1.333	1.338	-0.4	109	0.00
9		Benzaldehyde	0.911	0.840	7.8	113	0.00
10	C	Phenol	1.634	1.639	-0.3	110	0.00
11		bis(2-Chloroethyl)ether	1.318	1.335	-1.3	112	0.00
12		1,3-Dichlorobenzene	1.487	1.505	-1.2	110	0.00
13	C	1,4-Dichlorobenzene	1.514	1.521	-0.5	110	0.00
14		1,2-Dichlorobenzene	1.464	1.447	1.2	109	0.00
15		Benzyl Alcohol	1.069	1.076	-0.7	110	0.00
16		2,2'-oxybis(1-Chloropropane	1.716	1.731	-0.9	110	0.00
17		2-Methylphenol	1.061	1.050	1.0	108	0.00
18		Hexachloroethane	0.545	0.550	-0.9	110	0.00
19	P	n-Nitroso-di-n-propylamine	0.941	0.935	0.6	108	0.00
20		3+4-Methylphenols	1.434	1.419	1.0	109	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	110	0.00
22		Acetophenone	0.494	0.504	-2.0	108	0.00
23	S	Nitrobenzene-d5	0.374	0.386	-3.2	109	0.00
24		Nitrobenzene	0.351	0.359	-2.3	108	0.00
25		Isophorone	0.585	0.604	-3.2	110	0.00
26	C	2-Nitrophenol	0.158	0.167	-5.7	111	0.00
27		2,4-Dimethylphenol	0.250	0.255	-2.0	108	0.00
28		bis(2-Chloroethoxy)methane	0.407	0.423	-3.9	110	0.00
29	C	2,4-Dichlorophenol	0.277	0.286	-3.2	108	0.00
30		1,2,4-Trichlorobenzene	0.319	0.328	-2.8	109	0.00
31		Naphthalene	1.016	1.034	-1.8	108	0.00
32		Benzoic acid	0.189	0.196	-3.7	109	0.00
33		4-Chloroaniline	0.407	0.418	-2.7	109	0.00
34	C	Hexachlorobutadiene	0.183	0.187	-2.2	108	0.00
35		Caprolactam	0.085	0.086	-1.2	110	0.00
36	C	4-Chloro-3-methylphenol	0.284	0.294	-3.5	111	0.00
37		2-Methylnaphthalene	0.661	0.672	-1.7	108	0.00
38		1-Methylnaphthalene	0.643	0.656	-2.0	109	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	111	0.00
40		1,2,4,5-Tetrachlorobenzene	0.571	0.584	-2.3	108	0.00
41	P	Hexachlorocyclopentadiene	0.308	0.323	-4.9	109	0.00
42	S	2,4,6-Tribromophenol	0.224	0.235	-4.9	114	0.00
43	C	2,4,6-Trichlorophenol	0.378	0.393	-4.0	109	0.00
44		2,4,5-Trichlorophenol	0.398	0.417	-4.8	111	0.00
45	S	2-Fluorobiphenyl	1.415	1.444	-2.0	107	0.00
46		1,1'-Biphenyl	1.521	1.562	-2.7	109	0.00
47		2-Chloronaphthalene	1.169	1.198	-2.5	109	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.322	0.340	-5.6	112	0.00
49		Acenaphthylene	1.801	1.853	-2.9	109	0.00
50		Dimethylphthalate	1.316	1.370	-4.1	113	0.00
51		2,6-Dinitrotoluene	0.275	0.293	-6.5	112	0.00
52	C	Acenaphthene	1.079	1.103	-2.2	109	0.00
53		3-Nitroaniline	0.331	0.352	-6.3	113	0.00
54	P	2,4-Dinitrophenol	0.136	0.152	-11.8	125	0.00
55		Dibenzofuran	1.728	1.765	-2.1	110	0.00
56	P	4-Nitrophenol	0.260	0.280	-7.7	115	0.00
57		2,4-Dinitrotoluene	0.360	0.386	-7.2	114	0.00
58		Fluorene	1.357	1.397	-2.9	111	0.00
59		2,3,4,6-Tetrachlorophenol	0.320	0.338	-5.6	114	0.00
60		Diethylphthalate	1.304	1.359	-4.2	114	0.00
61		4-Chlorophenyl-phenylether	0.660	0.678	-2.7	111	0.00
62		4-Nitroaniline	0.339	0.363	-7.1	113	0.00
63		Azobenzene	1.339	1.394	-4.1	111	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	115	0.00
65		4,6-Dinitro-2-methylphenol	0.100	0.110	-10.0	121	0.00
66	c	n-Nitrosodiphenylamine	0.587	0.602	-2.6	112	0.00
67		4-Bromophenyl-phenylether	0.201	0.206	-2.5	112	0.00
68		Hexachlorobenzene	0.233	0.240	-3.0	113	0.00
69		Atrazine	0.154	0.168	-9.1	113	0.00
70	C	Pentachlorophenol	0.130	0.142	-9.2	118	0.00
71		Phenanthrene	1.052	1.074	-2.1	113	0.00
72		Anthracene	1.023	1.055	-3.1	113	0.00
73		Carbazole	1.037	1.074	-3.6	114	0.00
74		Di-n-butylphthalate	1.056	1.143	-8.2	120	0.00
75	C	Fluoranthene	1.144	1.193	-4.3	116	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	118	0.00
77		Benzidine	0.351	0.373	-6.3	118	0.00
78		Pyrene	1.308	1.330	-1.7	115	0.00
79	S	Terphenyl-d14	1.028	1.066	-3.7	116	0.00
80		Butylbenzylphthalate	0.467	0.503	-7.7	124	0.00
81		Benzo(a)anthracene	1.266	1.289	-1.8	115	0.00
82		3,3'-Dichlorobenzidine	0.297	0.315	-6.1	121	0.00
83		Chrysene	1.208	1.236	-2.3	116	0.00
84		Bis(2-ethylhexyl)phthalate	0.658	0.703	-6.8	122	0.00
85	c	Di-n-octyl phthalate	1.044	1.113	-6.6	122	0.00
86	I	Perylene-d12	1.000	1.000	0.0	118	0.00
87		Indeno(1,2,3-cd)pyrene	1.461	1.487	-1.8	115	0.00
88		Benzo(b)fluoranthene	1.191	1.212	-1.8	114	0.00
89		Benzo(k)fluoranthene	1.232	1.263	-2.5	118	0.00
90	C	Benzo(a)pyrene	1.013	1.041	-2.8	116	0.00
91		Dibenzo(a,h)anthracene	1.214	1.229	-1.2	114	0.00
92		Benzo(g,h,i)perylene	1.198	1.219	-1.8	115	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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