Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP091622\
Data File : BP011723.D
Acq On : 16 Sep 2022 09:45
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Sep 16 15:25:39 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP091422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 14 19:12:50 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	108	-0.01
2		1,4-Dioxane	0.484	0.477	1.4	110	0.00
3		Pyridine	1.395	1.431	-2.6	112	0.00
4		n-Nitrosodimethylamine	0.535	0.553	-3.4	113	0.00
5	S	2-Fluorophenol	1.099	1.111	-1.1	111	0.00
6		Aniline	1.804	1.839	-1.9	110	0.00
7	S	Phenol-d6	1.465	1.499	-2.3	110	0.00
8		2-Chlorophenol	1.303	1.311	-0.6	109	0.00
9		Benzaldehyde	0.938	0.929	1.0	112	-0.01
10	C	Phenol	1.647	1.676	-1.8	110	0.00
11		bis(2-Chloroethyl)ether	1.312	1.324	-0.9	110	0.00
12		1,3-Dichlorobenzene	1.445	1.421	1.7	108	-0.01
13	C	1,4-Dichlorobenzene	1.469	1.449	1.4	108	-0.01
14		1,2-Dichlorobenzene	1.401	1.377	1.7	108	0.00
15		Benzyl Alcohol	1.046	1.079	-3.2	111	-0.01
16		2,2'-oxybis(1-Chloropropane	1.753	1.791	-2.2	111	-0.01
17		2-Methylphenol	1.070	1.090	-1.9	110	-0.01
18		Hexachloroethane	0.508	0.519	-2.2	111	0.00
19	P	n-Nitroso-di-n-propylamine	0.948	0.970	-2.3	108	0.00
20		3+4-Methylphenols	1.469	1.490	-1.4	109	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	109	-0.01
22		Acetophenone	0.469	0.471	-0.4	109	0.00
23	S	Nitrobenzene-d5	0.337	0.346	-2.7	110	0.00
24		Nitrobenzene	0.351	0.358	-2.0	110	0.00
25		Isophorone	0.597	0.605	-1.3	108	0.00
26	C	2-Nitrophenol	0.144	0.144	0.0	105	0.00
27		2,4-Dimethylphenol	0.249	0.254	-2.0	108	0.00
28		bis(2-Chloroethoxy)methane	0.381	0.384	-0.8	109	0.00
29	C	2,4-Dichlorophenol	0.276	0.279	-1.1	108	0.00
30		1,2,4-Trichlorobenzene	0.306	0.299	2.3	107	0.00
31		Naphthalene	1.038	1.024	1.3	108	0.00
32		Benzoic acid	0.187	0.177	5.3	102	0.00
33		4-Chloroaniline	0.407	0.409	-0.5	108	0.00
34	C	Hexachlorobutadiene	0.168	0.162	3.6	106	0.00
35		Caprolactam	0.090	0.087	3.3	104	0.00
36	C	4-Chloro-3-methylphenol	0.297	0.299	-0.7	107	0.00
37		2-Methylnaphthalene	0.695	0.689	0.9	108	0.00
38		1-Methylnaphthalene	0.679	0.672	1.0	108	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	107	0.00
40		1,2,4,5-Tetrachlorobenzene	0.528	0.520	1.5	106	0.00
41	P	Hexachlorocyclopentadiene	0.263	0.262	0.4	105	0.00
42	S	2,4,6-Tribromophenol	0.156	0.153	1.9	102	0.00
43	C	2,4,6-Trichlorophenol	0.359	0.365	-1.7	106	0.00
44		2,4,5-Trichlorophenol	0.401	0.404	-0.7	106	0.00
45	S	2-Fluorobiphenyl	1.287	1.301	-1.1	108	0.00
46		1,1'-Biphenyl	1.455	1.464	-0.6	108	-0.01
47		2-Chloronaphthalene	1.145	1.143	0.2	107	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.321	0.345	-7.5	109	0.00
49		Acenaphthylene	1.771	1.818	-2.7	108	0.00
50		Dimethylphthalate	1.395	1.405	-0.7	108	0.00
51		2,6-Dinitrotoluene	0.283	0.293	-3.5	106	0.00
52	C	Acenaphthene	1.178	1.188	-0.8	107	-0.01
53		3-Nitroaniline	0.326	0.346	-6.1	108	0.00
54	P	2,4-Dinitrophenol	0.138	0.128	7.2	100	0.00
55		Dibenzofuran	1.697	1.697	0.0	108	0.00
56	P	4-Nitrophenol	0.275	0.279	-1.5	106	0.00
57		2,4-Dinitrotoluene	0.366	0.392	-7.1	107	0.00
58		Fluorene	1.376	1.408	-2.3	108	0.00
59		2,3,4,6-Tetrachlorophenol	0.330	0.326	1.2	104	0.00
60		Diethylphthalate	1.389	1.425	-2.6	109	0.00
61		4-Chlorophenyl-phenylether	0.636	0.640	-0.6	108	0.00
62		4-Nitroaniline	0.328	0.361	-10.1	110	0.00
63		Azobenzene	1.335	1.421	-6.4	110	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	107	0.00
65		4,6-Dinitro-2-methylphenol	0.098	0.094	4.1	103	0.00
66	c	n-Nitrosodiphenylamine	0.595	0.601	-1.0	107	0.00
67		4-Bromophenyl-phenylether	0.187	0.181	3.2	104	0.00
68		Hexachlorobenzene	0.202	0.193	4.5	104	0.00
69		Atrazine	0.182	0.181	0.5	104	0.00
70	C	Pentachlorophenol	0.128	0.122	4.7	102	0.00
71		Phenanthrene	1.088	1.083	0.5	107	0.00
72		Anthracene	1.029	1.047	-1.7	107	0.00
73		Carbazole	0.971	0.992	-2.2	107	0.00
74		Di-n-butylphthalate	1.170	1.219	-4.2	107	0.00
75	C	Fluoranthene	1.174	1.187	-1.1	107	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	106	0.00
77		Benzidine	0.476	0.557	-17.0	107	0.00
78		Pyrene	1.349	1.381	-2.4	108	0.00
79	S	Terphenyl-d14	0.920	0.938	-2.0	107	0.00
80		Butylbenzylphthalate	0.565	0.601	-6.4	107	0.00
81		Benzo(a)anthracene	1.289	1.305	-1.2	107	0.00
82		3,3'-Dichlorobenzidine	0.392	0.408	-4.1	105	0.00
83		Chrysene	1.232	1.234	-0.2	106	0.00
84		Bis(2-ethylhexyl)phthalate	0.813	0.882	-8.5	107	0.00
85	c	Di-n-octyl phthalate	1.395	1.421	-1.9	105	0.00
86	I	Perylene-d12	1.000	1.000	0.0	104	0.00
87		Indeno(1,2,3-cd)pyrene	1.441	1.476	-2.4	105	0.00
88		Benzo(b)fluoranthene	1.237	1.295	-4.7	107	0.00
89		Benzo(k)fluoranthene	1.247	1.265	-1.4	104	0.00
90	C	Benzo(a)pyrene	1.022	1.058	-3.5	105	0.00
91		Dibenzo(a,h)anthracene	1.196	1.229	-2.8	105	0.00
92		Benzo(g,h,i)perylene	1.166	1.176	-0.9	104	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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