Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111121\
Data File : BP007942.D
Acq On : 11 Nov 2021 13:07
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Nov 11 13:54:36 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Nov 11 13:53:53 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	59	0.00
2		1,4-Dioxane	0.447	0.438	2.0	63	0.00
3		Pyridine	1.308	1.372	-4.9	66	0.00
4		n-Nitrosodimethylamine	0.560	0.548	2.1	61	0.00
5	S	2-Fluorophenol	1.179	1.143	3.1	59	0.00
6		Aniline	1.923	1.899	1.2	59	0.00
7	S	Phenol-d6	1.707	1.673	2.0	59	0.00
8		2-Chlorophenol	1.300	1.280	1.5	60	0.00
9		Benzaldehyde	0.928	0.862	7.1	58	0.00
10	C	Phenol	1.630	1.607	1.4	59	0.00
11		bis(2-Chloroethyl)ether	1.301	1.239	4.8	57	0.00
12		1,3-Dichlorobenzene	1.445	1.400	3.1	59	0.00
13	C	1,4-Dichlorobenzene	1.477	1.442	2.4	60	0.00
14		1,2-Dichlorobenzene	1.427	1.410	1.2	60	0.00
15		Benzyl Alcohol	1.292	1.318	-2.0	59	0.00
16		2,2'-oxybis(1-Chloropropane	1.525	1.429	6.3	56	0.00
17		2-Methylphenol	1.127	1.126	0.1	59	0.00
18		Hexachloroethane	0.511	0.494	3.3	58	0.00
19	P	n-Nitroso-di-n-propylamine	1.045	1.061	-1.5	60	0.00
20		3+4-Methylphenols	1.586	1.620	-2.1	60	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	59	0.00
22		Acetophenone	0.504	0.504	0.0	60	0.00
23	S	Nitrobenzene-d5	0.372	0.374	-0.5	59	0.00
24		Nitrobenzene	0.354	0.355	-0.3	60	0.00
25		Isophorone	0.677	0.682	-0.7	59	0.00
26	C	2-Nitrophenol	0.179	0.188	-5.0	60	0.00
27		2,4-Dimethylphenol	0.276	0.276	0.0	59	0.00
28		bis(2-Chloroethoxy)methane	0.446	0.424	4.9	56	0.00
29	C	2,4-Dichlorophenol	0.332	0.345	-3.9	60	0.00
30		1,2,4-Trichlorobenzene	0.369	0.370	-0.3	60	0.00
31		Naphthalene	1.012	1.003	0.9	60	0.00
32		Benzoic acid	0.238	0.267	-12.2	59	0.00
33		4-Chloroaniline	0.446	0.457	-2.5	61	0.00
34	C	Hexachlorobutadiene	0.242	0.256	-5.8	63	0.00
35		Caprolactam	0.116	0.125	-7.8	62	0.00
36	C	4-Chloro-3-methylphenol	0.375	0.398	-6.1	61	0.00
37		2-Methylnaphthalene	0.746	0.756	-1.3	60	0.00
38		1-Methylnaphthalene	0.731	0.739	-1.1	60	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	62	0.00
40		1,2,4,5-Tetrachlorobenzene	0.627	0.637	-1.6	64	0.00
41	P	Hexachlorocyclopentadiene	0.336	0.309	8.0	54	0.00
42	S	2,4,6-Tribromophenol	0.285	0.315	-10.5	69	0.00
43	C	2,4,6-Trichlorophenol	0.438	0.453	-3.4	63	0.00
44		2,4,5-Trichlorophenol	0.474	0.497	-4.9	64	0.00
45	S	2-Fluorobiphenyl	1.373	1.356	1.2	62	0.00
46		1,1'-Biphenyl	1.444	1.425	1.3	62	0.00
47		2-Chloronaphthalene	1.122	1.117	0.4	62	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.323	0.336	-4.0	62	0.00
49		Acenaphthylene	1.762	1.762	0.0	61	0.00
50		Dimethylphthalate	1.544	1.547	-0.2	62	0.00
51		2,6-Dinitrotoluene	0.319	0.330	-3.4	62	0.00
52	C	Acenaphthene	1.064	1.050	1.3	62	0.00
53		3-Nitroaniline	0.341	0.346	-1.5	61	0.00
54	P	2,4-Dinitrophenol	0.196	0.217	-10.7	64	0.00
55		Dibenzofuran	1.800	1.662	7.7	58	0.00
56	P	4-Nitrophenol	0.289	0.285	1.4	59	0.00
57		2,4-Dinitrotoluene	0.439	0.434	1.1	59	0.00
58		Fluorene	1.366	1.295	5.2	58	0.00
59		2,3,4,6-Tetrachlorophenol	0.427	0.431	-0.9	61	0.00
60		Diethylphthalate	1.499	1.444	3.7	59	0.00
61		4-Chlorophenyl-phenylether	0.746	0.729	2.3	60	0.00
62		4-Nitroaniline	0.346	0.340	1.7	58	0.00
63		Azobenzene	1.265	1.210	4.3	60	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	62	0.00
65		4,6-Dinitro-2-methylphenol	0.131	0.140	-6.9	60	0.00
66	c	n-Nitrosodiphenylamine	0.569	0.565	0.7	62	0.00
67		4-Bromophenyl-phenylether	0.221	0.233	-5.4	66	0.00
68		Hexachlorobenzene	0.247	0.265	-7.3	67	0.00
69		Atrazine	0.218	0.231	-6.0	65	0.00
70	C	Pentachlorophenol	0.159	0.171	-7.5	63	0.00
71		Phenanthrene	1.051	1.042	0.9	62	0.00
72		Anthracene	1.035	1.042	-0.7	63	0.00
73		Carbazole	1.029	1.015	1.4	62	0.00
74		Di-n-butylphthalate	1.171	1.188	-1.5	63	0.00
75	C	Fluoranthene	1.326	1.337	-0.8	64	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	69	0.00
77		Benzidine	0.504	0.498	1.2	61	0.00
78		Pyrene	1.227	1.154	5.9	65	0.00
79	S	Terphenyl-d14	0.982	0.963	1.9	68	0.00
80		Butylbenzylphthalate	0.516	0.497	3.7	65	0.00
81		Benzo(a)anthracene	1.318	1.280	2.9	68	0.00
82		3,3'-Dichlorobenzidine	0.442	0.450	-1.8	70	0.00
83		Chrysene	1.247	1.225	1.8	69	0.00
84		Bis(2-ethylhexyl)phthalate	0.739	0.712	3.7	66	0.00
85	c	Di-n-octyl phthalate	1.330	1.332	-0.2	68	0.00
86	I	Perylene-d12	1.000	1.000	0.0	70	0.00
87		Indeno(1,2,3-cd)pyrene	1.491	1.442	3.3	69	0.00
88		Benzo(b)fluoranthene	1.179	1.153	2.2	69	0.00
89		Benzo(k)fluoranthene	1.170	1.166	0.3	71	0.00
90	C	Benzo(a)pyrene	1.021	0.999	2.2	69	0.00
91		Dibenzo(a,h)anthracene	1.236	1.196	3.2	69	0.00
92		Benzo(g,h,i)perylene	1.226	1.183	3.5	69	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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