Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111021\
Data File : BP007912.D
Acq On : 10 Nov 2021 10:28
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Nov 10 13:12:26 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP110221.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Nov 09 10:45:24 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	72	0.01
2		1,4-Dioxane	0.447	0.437	2.2	77	0.00
3		Pyridine	1.308	1.346	-2.9	79	0.01
4		n-Nitrosodimethylamine	0.560	0.546	2.5	74	0.00
5	S	2-Fluorophenol	1.179	1.151	2.4	73	0.01
6		Aniline	1.923	1.831	4.8	69	0.00
7	S	Phenol-d6	1.707	1.614	5.4	69	0.01
8		2-Chlorophenol	1.300	1.272	2.2	72	0.01
9		Benzaldehyde	0.928	0.829	10.7	68	0.00
10	C	Phenol	1.630	1.542	5.4	69	0.00
11		bis(2-Chloroethyl)ether	1.301	1.210	7.0	68	0.00
12		1,3-Dichlorobenzene	1.445	1.412	2.3	73	0.01
13	C	1,4-Dichlorobenzene	1.477	1.433	3.0	73	0.01
14		1,2-Dichlorobenzene	1.427	1.370	4.0	71	0.01
15		Benzyl Alcohol	1.292	1.221	5.5	67	0.00
16		2,2'-oxybis(1-Chloropropane	1.525	1.311	14.0	63	0.00
17		2-Methylphenol	1.127	1.063	5.7	68	0.00
18		Hexachloroethane	0.511	0.463	9.4	66	0.00
19	P	n-Nitroso-di-n-propylamine	1.045	0.890	14.8	61	0.01
20		3+4-Methylphenols	1.586	1.398	11.9	63	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	62	0.01
22		Acetophenone	0.504	0.503	0.2	63	0.01
23	S	Nitrobenzene-d5	0.372	0.380	-2.2	64	0.01
24		Nitrobenzene	0.354	0.360	-1.7	64	0.01
25		Isophorone	0.677	0.659	2.7	60	0.01
26	C	2-Nitrophenol	0.179	0.189	-5.6	64	0.01
27		2,4-Dimethylphenol	0.276	0.273	1.1	62	0.00
28		bis(2-Chloroethoxy)methane	0.446	0.410	8.1	58	0.01
29	C	2,4-Dichlorophenol	0.332	0.335	-0.9	62	0.01
30		1,2,4-Trichlorobenzene	0.369	0.372	-0.8	64	0.01
31		Naphthalene	1.012	1.000	1.2	63	0.00
32		Benzoic acid	0.238	0.247	-3.8	58	0.01
33		4-Chloroaniline	0.446	0.441	1.1	62	0.01
34	C	Hexachlorobutadiene	0.242	0.256	-5.8	67	0.01
35		Caprolactam	0.116	0.116	0.0	61	0.00
36	C	4-Chloro-3-methylphenol	0.375	0.371	1.1	60	0.01
37		2-Methylnaphthalene	0.746	0.726	2.7	61	0.00
38		1-Methylnaphthalene	0.731	0.712	2.6	62	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	61	0.00
40		1,2,4,5-Tetrachlorobenzene	0.627	0.646	-3.0	64	0.01
41	P	Hexachlorocyclopentadiene	0.336	0.308	8.3	53	0.00
42	S	2,4,6-Tribromophenol	0.285	0.328	-15.1	71	0.00
43	C	2,4,6-Trichlorophenol	0.438	0.457	-4.3	62	0.00
44		2,4,5-Trichlorophenol	0.474	0.490	-3.4	62	0.01
45	S	2-Fluorobiphenyl	1.373	1.387	-1.0	63	0.00
46		1,1'-Biphenyl	1.444	1.443	0.1	62	0.00
47		2-Chloronaphthalene	1.122	1.139	-1.5	62	0.01

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.323	0.343	-6.2	63	0.01
49		Acenaphthylene	1.762	1.792	-1.7	62	0.01
50		Dimethylphthalate	1.544	1.547	-0.2	62	0.00
51		2,6-Dinitrotoluene	0.319	0.329	-3.1	61	0.00
52	C	Acenaphthene	1.064	1.065	-0.1	62	0.00
53		3-Nitroaniline	0.341	0.347	-1.8	61	0.00
54	P	2,4-Dinitrophenol	0.196	0.210	-7.1	61	0.01
55		Dibenzofuran	1.800	1.805	-0.3	62	0.00
56	P	4-Nitrophenol	0.289	0.291	-0.7	60	0.00
57		2,4-Dinitrotoluene	0.439	0.466	-6.2	62	0.01
58		Fluorene	1.366	1.399	-2.4	62	0.01
59		2,3,4,6-Tetrachlorophenol	0.427	0.458	-7.3	64	0.00
60		Diethylphthalate	1.499	1.551	-3.5	62	0.00
61		4-Chlorophenyl-phenylether	0.746	0.768	-2.9	63	0.01
62		4-Nitroaniline	0.346	0.372	-7.5	63	0.00
63		Azobenzene	1.265	1.297	-2.5	63	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	65	0.00
65		4,6-Dinitro-2-methylphenol	0.131	0.139	-6.1	62	0.00
66	c	n-Nitrosodiphenylamine	0.569	0.561	1.4	65	0.00
67		4-Bromophenyl-phenylether	0.221	0.230	-4.1	68	0.00
68		Hexachlorobenzene	0.247	0.256	-3.6	69	0.01
69		Atrazine	0.218	0.231	-6.0	69	0.00
70	C	Pentachlorophenol	0.159	0.168	-5.7	65	0.00
71		Phenanthrene	1.051	1.042	0.9	66	0.00
72		Anthracene	1.035	1.045	-1.0	67	0.00
73		Carbazole	1.029	1.022	0.7	67	0.00
74		Di-n-butylphthalate	1.171	1.189	-1.5	67	0.00
75	C	Fluoranthene	1.326	1.346	-1.5	68	0.01
76	I	Chrysene-d12	1.000	1.000	0.0	72	0.00
77		Benzidine	0.504	0.491	2.6	63	0.00
78		Pyrene	1.227	1.172	4.5	69	0.00
79	S	Terphenyl-d14	0.982	0.950	3.3	71	0.00
80		Butylbenzylphthalate	0.516	0.509	1.4	70	0.00
81		Benzo(a)anthracene	1.318	1.295	1.7	72	0.00
82		3,3'-Dichlorobenzidine	0.442	0.431	2.5	70	0.00
83		Chrysene	1.247	1.225	1.8	72	0.00
84		Bis(2-ethylhexyl)phthalate	0.739	0.724	2.0	70	0.00
85	c	Di-n-octyl phthalate	1.330	1.333	-0.2	71	0.00
86	I	Perylene-d12	1.000	1.000	0.0	74	0.00
87		Indeno(1,2,3-cd)pyrene	1.491	1.458	2.2	73	0.02
88		Benzo(b)fluoranthene	1.179	1.175	0.3	74	0.01
89		Benzo(k)fluoranthene	1.170	1.132	3.2	73	0.00
90	C	Benzo(a)pyrene	1.021	1.004	1.7	73	0.01
91		Dibenzo(a,h)anthracene	1.236	1.207	2.3	74	0.02
92		Benzo(g,h,i)perylene	1.226	1.198	2.3	73	0.02

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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