Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP042121\
Data File : BP005363.D
Acq On : 21 Apr 2021 17:30
Operator : CG/JU
Sample : SSTDICV040
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_P
ClientSampleId : ICVBP042121

Quant Time: Apr 21 18:37:41 2021
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP042121.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Apr 21 17:40:15 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	94	0.00
2		1,4-Dioxane	0.614	0.685	-11.6	102	0.00
3		Pyridine	1.332	1.656	-24.3	110	0.00
4		n-Nitrosodimethylamine	0.686	0.826	-20.4	115	0.00
5	S	2-Fluorophenol	1.151	1.204	-4.6	98	0.00
6		Aniline	1.721	1.847	-7.3	99	0.00
7	S	Phenol-d6	1.586	1.679	-5.9	99	0.00
8		2-Chlorophenol	1.179	1.194	-1.3	96	0.00
9		Benzaldehyde	0.928	1.015	-9.4	100	0.00
10	C	Phenol	1.582	1.674	-5.8	99	0.00
11		bis(2-Chloroethyl)ether	1.141	1.203	-5.4	98	0.00
12		1,3-Dichlorobenzene	1.425	1.413	0.8	94	0.01
13	C	1,4-Dichlorobenzene	1.454	1.424	2.1	94	0.00
14		1,2-Dichlorobenzene	1.390	1.400	-0.7	96	0.00
15		Benzyl Alcohol	1.265	1.468	-16.0	106	0.01
16		2,2'-oxybis(1-Chloropropane	0.957	1.037	-8.4	102	0.00
17		2-Methylphenol	0.995	1.058	-6.3	97	0.00
18		Hexachloroethane	0.556	0.583	-4.9	97	0.00
19	P	n-Nitroso-di-n-propylamine	1.030	1.216	-18.1	105	0.00
20		3+4-Methylphenols	1.336	1.440	-7.8	98	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	94	0.00
22		Acetophenone	0.574	0.618	-7.7	99	0.01
23	S	Nitrobenzene-d5	0.460	0.544	-18.3	101	0.01
24		Nitrobenzene	0.497	0.568	-14.3	101	0.01
25		Isophorone	0.799	0.927	-16.0	101	0.00
26	C	2-Nitrophenol	0.160	0.188	-17.5	102	0.00
27		2,4-Dimethylphenol	0.270	0.282	-4.4	95	0.01
28		bis(2-Chloroethoxy)methane	0.384	0.414	-7.8	97	0.00
29	C	2,4-Dichlorophenol	0.357	0.387	-8.4	98	0.00
30		1,2,4-Trichlorobenzene	0.469	0.474	-1.1	94	0.00
31		Naphthalene	1.055	1.071	-1.5	94	0.00
32		Benzoic acid	0.176	0.201	-14.2	104	0.02
33		4-Chloroaniline	0.400	0.437	-9.2	98	0.01
34	C	Hexachlorobutadiene	0.416	0.424	-1.9	94	0.01
35		Caprolactam	0.100	0.106	-6.0	96	0.01
36	C	4-Chloro-3-methylphenol	0.385	0.425	-10.4	97	0.00
37		2-Methylnaphthalene	0.773	0.795	-2.8	94	0.00
38		1-Methylnaphthalene	0.738	0.760	-3.0	95	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	93	0.01
40		1,2,4,5-Tetrachlorobenzene	0.871	0.918	-5.4	95	0.00
41	P	Hexachlorocyclopentadiene	0.282	0.367	-30.1#	108	0.00
42	S	2,4,6-Tribromophenol	0.383	0.414	-8.1	94	0.01
43	C	2,4,6-Trichlorophenol	0.472	0.539	-14.2	98	0.01
44		2,4,5-Trichlorophenol	0.515	0.582	-13.0	99	0.00
45	S	2-Fluorobiphenyl	1.545	1.651	-6.9	96	0.01
46		1,1'-Biphenyl	1.423	1.515	-6.5	95	0.00
47		2-Chloronaphthalene	1.172	1.254	-7.0	95	0.00

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		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.322	0.416	-29.2#	105	0.00
49		Acenaphthylene	1.696	1.856	-9.4	94	0.00
50		Dimethylphthalate	1.452	1.571	-8.2	95	0.00
51		2,6-Dinitrotoluene	0.317	0.364	-14.8	97	0.01
52	C	Acenaphthene	1.070	1.126	-5.2	94	0.01
53		3-Nitroaniline	0.262	0.299	-14.1	94	0.01
54	P	2,4-Dinitrophenol	0.180	0.208	-15.6	100	0.01
55		Dibenzofuran	1.917	1.988	-3.7	93	0.00
56	P	4-Nitrophenol	0.208	0.229	-10.1	96	0.01
57		2,4-Dinitrotoluene	0.443	0.501	-13.1	95	0.01
58		Fluorene	1.538	1.641	-6.7	95	0.00
59		2,3,4,6-Tetrachlorophenol	0.536	0.598	-11.6	95	0.00
60		Diethylphthalate	1.370	1.543	-12.6	98	0.00
61		4-Chlorophenyl-phenylether	0.987	1.054	-6.8	95	0.01
62		4-Nitroaniline	0.246	0.280	-13.8	91	0.00
63		Azobenzene	1.597	1.871	-17.2	99	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	93	0.00
65		4,6-Dinitro-2-methylphenol	0.131	0.145	-10.7	97	0.00
66	c	n-Nitrosodiphenylamine	0.547	0.571	-4.4	93	0.00
67		4-Bromophenyl-phenylether	0.273	0.288	-5.5	94	0.01
68		Hexachlorobenzene	0.320	0.327	-2.2	93	0.01
69		Atrazine	0.228	0.259	-13.6	96	0.00
70	C	Pentachlorophenol	0.172	0.183	-6.4	96	0.01
71		Phenanthrene	1.066	1.081	-1.4	92	0.01
72		Anthracene	1.038	1.072	-3.3	93	0.00
73		Carbazole	0.934	0.958	-2.6	92	0.01
74		Di-n-butylphthalate	0.882	1.005	-13.9	98	0.01
75	C	Fluoranthene	1.463	1.492	-2.0	91	0.01
76	I	Chrysene-d12	1.000	1.000	0.0	90	0.00
77		Benzidine	0.316	0.369	-16.8	97	0.00
78		Pyrene	1.121	1.164	-3.8	92	0.01
79	S	Terphenyl-d14	1.079	1.117	-3.5	92	0.00
80		Butylbenzylphthalate	0.255	0.299	-17.3	99	0.00
81		Benzo(a)anthracene	1.287	1.297	-0.8	90	0.00
82		3,3'-Dichlorobenzidine	0.400	0.427	-6.7	92	0.01
83		Chrysene	1.282	1.275	0.5	90	0.00
84		Bis(2-ethylhexyl)phthalate	0.391	0.476	-21.7	100	0.00
85	c	Di-n-octyl phthalate	0.760	0.820	-7.9	98	0.00
86	I	Perylene-d12	1.000	1.000	0.0	90	0.01
87		Indeno(1,2,3-cd)pyrene	1.528	1.536	-0.5	89	0.02
88		Benzo(b)fluoranthene	1.317	1.350	-2.5	90	0.01
89		Benzo(k)fluoranthene	1.293	1.297	-0.3	88	0.01
90	C	Benzo(a)pyrene	1.201	1.216	-1.2	89	0.01
91		Dibenzo(a,h)anthracene	1.323	1.332	-0.7	90	0.02
92		Benzo(g,h,i)perylene	1.239	1.249	-0.8	89	0.02

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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