Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP011822\
Data File : BP008834.D
Acq On : 18 Jan 2022 11:23
Operator : CG/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Jan 19 03:35:06 2022
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP011422.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Jan 14 18:05:00 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	99	0.00
2		1,4-Dioxane	0.523	0.502	4.0	112	0.00
3		Pyridine	1.096	1.187	-8.3	119	0.00
4		n-Nitrosodimethylamine	0.514	0.555	-8.0	119	0.00
5	S	2-Fluorophenol	1.293	1.311	-1.4	114	0.00
6		Aniline	1.955	2.128	-8.8	119	0.00
7	S	Phenol-d6	1.566	1.699	-8.5	118	0.00
8		2-Chlorophenol	1.377	1.443	-4.8	115	0.00
9		Benzaldehyde	1.047	1.071	-2.3	113	0.00
10	C	Phenol	1.597	1.728	-8.2	118	0.00
11		bis(2-Chloroethyl)ether	1.270	1.349	-6.2	117	0.00
12		1,3-Dichlorobenzene	1.640	1.648	-0.5	112	0.00
13	C	1,4-Dichlorobenzene	1.662	1.684	-1.3	113	0.00
14		1,2-Dichlorobenzene	1.585	1.618	-2.1	114	0.00
15		Benzyl Alcohol	1.099	1.263	-14.9	123	0.00
16		2,2'-oxybis(1-Chloropropane	1.421	1.538	-8.2	118	0.00
17		2-Methylphenol	1.070	1.182	-10.5	120	0.00
18		Hexachloroethane	0.564	0.582	-3.2	113	0.00
19	P	n-Nitroso-di-n-propylamine	0.915	1.072	-17.2	124	0.00
20		3+4-Methylphenols	1.420	1.618	-13.9	122	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	106	0.00
22		Acetophenone	0.509	0.528	-3.7	122	0.00
23	S	Nitrobenzene-d5	0.352	0.367	-4.3	122	0.00
24		Nitrobenzene	0.356	0.368	-3.4	122	0.00
25		Isophorone	0.637	0.703	-10.4	128	0.00
26	C	2-Nitrophenol	0.187	0.200	-7.0	123	0.00
27		2,4-Dimethylphenol	0.319	0.337	-5.6	123	0.00
28		bis(2-Chloroethoxy)methane	0.415	0.438	-5.5	124	0.00
29	C	2,4-Dichlorophenol	0.348	0.365	-4.9	122	0.00
30		1,2,4-Trichlorobenzene	0.408	0.401	1.7	118	0.00
31		Naphthalene	1.086	1.090	-0.4	120	0.00
32		Benzoic acid	0.183	0.223	-21.9	133	0.01
33		4-Chloroaniline	0.443	0.479	-8.1	126	0.00
34	C	Hexachlorobutadiene	0.255	0.248	2.7	116	0.00
35		Caprolactam	0.096	0.121	-26.0#	145	0.01
36	C	4-Chloro-3-methylphenol	0.314	0.356	-13.4	131	0.00
37		2-Methylnaphthalene	0.794	0.842	-6.0	125	0.00
38		1-Methylnaphthalene	0.729	0.778	-6.7	125	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	117	0.00
40		1,2,4,5-Tetrachlorobenzene	0.729	0.688	5.6	124	0.00
41	P	Hexachlorocyclopentadiene	0.373	0.347	7.0	113	0.00
42	S	2,4,6-Tribromophenol	0.291	0.321	-10.3	144	0.00
43	C	2,4,6-Trichlorophenol	0.451	0.459	-1.8	131	0.00
44		2,4,5-Trichlorophenol	0.485	0.497	-2.5	132	0.00
45	S	2-Fluorobiphenyl	1.517	1.446	4.7	127	0.00
46		1,1'-Biphenyl	1.627	1.576	3.1	128	0.00
47		2-Chloronaphthalene	1.255	1.217	3.0	128	0.00

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Max. RRF Dev : 25% Max. Rel. Area : 150%

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.279	0.312	-11.8	143	0.00
49		Acenaphthylene	1.896	1.910	-0.7	132	0.00
50		Dimethylphthalate	1.485	1.554	-4.6	138	0.00
51		2,6-Dinitrotoluene	0.315	0.342	-8.6	140	0.00
52	C	Acenaphthene	1.124	1.139	-1.3	133	0.00
53		3-Nitroaniline	0.310	0.351	-13.2	146	0.00
54	P	2,4-Dinitrophenol	0.127	0.160	-26.0#	156#	0.00
55		Dibenzofuran	1.832	1.866	-1.9	134	0.00
56	P	4-Nitrophenol	0.223	0.269	-20.6	152#	0.00
57		2,4-Dinitrotoluene	0.405	0.477	-17.8	149	0.00
58		Fluorene	1.447	1.520	-5.0	137	0.00
59		2,3,4,6-Tetrachlorophenol	0.400	0.433	-8.2	140	0.00
60		Diethylphthalate	1.409	1.534	-8.9	143	0.00
61		4-Chlorophenyl-phenylether	0.788	0.826	-4.8	136	0.00
62		4-Nitroaniline	0.300	0.360	-20.0	154#	0.00
63		Azobenzene	1.170	1.294	-10.6	142	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	132	0.00
65		4,6-Dinitro-2-methylphenol	0.100	0.114	-14.0	157#	0.00
66	c	n-Nitrosodiphenylamine	0.644	0.636	1.2	144	0.00
67		4-Bromophenyl-phenylether	0.250	0.245	2.0	141	0.00
68		Hexachlorobenzene	0.286	0.277	3.1	141	0.00
69		Atrazine	0.239	0.255	-6.7	155#	0.00
70	C	Pentachlorophenol	0.152	0.163	-7.2	153#	0.00
71		Phenanthrene	1.126	1.128	-0.2	148	0.00
72		Anthracene	1.130	1.167	-3.3	151#	0.00
73		Carbazole	0.971	1.035	-6.6	158#	0.00
74		Di-n-butylphthalate	1.169	1.258	-7.6	155#	0.00
75	C	Fluoranthene	1.252	1.357	-8.4	164#	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	152#	0.00
77		Benzidine	0.707	0.792	-12.0	170#	0.00
78		Pyrene	1.396	1.416	-1.4	166#	0.00
79	S	Terphenyl-d14	1.185	1.141	3.7	158#	0.00
80		Butylbenzylphthalate	0.561	0.594	-5.9	174#	0.00
81		Benzo(a)anthracene	1.346	1.358	-0.9	172#	0.00
82		3,3'-Dichlorobenzidine	0.581	0.605	-4.1	174#	0.00
83		Chrysene	1.304	1.304	0.0	170#	0.00
84		Bis(2-ethylhexyl)phthalate	0.869	0.887	-2.1	167#	0.00
85	c	Di-n-octyl phthalate	1.470	1.491	-1.4	183#	0.00
86	I	Perylene-d12	1.000	1.000	0.0	137	0.00
87		Indeno(1,2,3-cd)pyrene	1.648	1.407	14.6	132	0.00
88		Benzo(b)fluoranthene	1.336	1.456	-9.0	169#	0.00
89		Benzo(k)fluoranthene	1.293	1.377	-6.5	170#	0.00
90	C	Benzo(a)pyrene	1.289	1.341	-4.0	162#	0.00
91		Dibenzo(a,h)anthracene	1.401	1.191	15.0	131	0.00
92		Benzo(g,h,i)perylene	1.368	1.098	19.7	123	0.00

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Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)

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