Evaluate Continuing Calibration Report

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020624\
Data File : BP019582.D
Acq On : 06 Feb 2024 19:50
Operator : MA/JU
Sample : SSTDCCC040
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument :
BNA_P
LabSampleId :
SSTDCCC040

Quant Time: Feb 07 01:08:39 2024
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\8270E-BP013124.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Feb 01 00:51:48 2024
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	120	0.00
2		1,4-Dioxane	0.478	0.440	7.9	126	0.00
3		Pyridine	1.296	1.099	15.2	110	0.00
4		n-Nitrosodimethylamine	0.564	0.588	-4.3	133	0.00
5	S	2-Fluorophenol	1.241	1.142	8.0	121	0.00
6		Aniline	1.626	1.577	3.0	122	0.00
7	S	Phenol-d6	1.673	1.631	2.5	124	0.00
8		2-Chlorophenol	1.274	1.241	2.6	125	0.00
9		Benzaldehyde	1.173	1.021	13.0	92	0.00
10	C	Phenol	1.667	1.613	3.2	123	0.00
11		bis(2-Chloroethyl)ether	1.297	1.298	-0.1	125	0.00
12		1,3-Dichlorobenzene	1.559	1.507	3.3	125	-0.01
13	C	1,4-Dichlorobenzene	1.579	1.530	3.1	126	0.00
14		1,2-Dichlorobenzene	1.528	1.494	2.2	126	0.00
15		Benzyl Alcohol	1.337	1.304	2.5	120	0.00
16		2,2'-oxybis(1-Chloropropane	1.297	1.286	0.8	121	0.00
17		2-Methylphenol	1.123	1.096	2.4	120	0.00
18		Hexachloroethane	0.612	0.606	1.0	126	0.00
19	P	n-Nitroso-di-n-propylamine	1.143	1.099	3.8	117	0.00
20		3+4-Methylphenols	1.532	1.511	1.4	123	0.00
21	I	Naphthalene-d8	1.000	1.000	0.0	120	0.00
22		Acetophenone	0.570	0.546	4.2	120	0.00
23	S	Nitrobenzene-d5	0.462	0.447	3.2	122	0.00
24		Nitrobenzene	0.465	0.452	2.8	123	0.00
25		Isophorone	0.804	0.767	4.6	118	0.00
26	C	2-Nitrophenol	0.177	0.179	-1.1	123	0.00
27		2,4-Dimethylphenol	0.227	0.213	6.2	117	-0.01
28		bis(2-Chloroethoxy)methane	0.455	0.436	4.2	118	0.00
29	C	2,4-Dichlorophenol	0.343	0.332	3.2	121	0.00
30		1,2,4-Trichlorobenzene	0.420	0.412	1.9	125	-0.01
31		Naphthalene	1.097	1.060	3.4	123	-0.01
32		Benzoic acid	0.221	0.227	-2.7	123	0.00
33		4-Chloroaniline	0.406	0.391	3.7	120	0.00
34	C	Hexachlorobutadiene	0.311	0.314	-1.0	127	0.00
35		Caprolactam	0.097	0.099	-2.1	136	0.00
36	C	4-Chloro-3-methylphenol	0.373	0.368	1.3	124	0.00
37		2-Methylnaphthalene	0.759	0.735	3.2	121	0.00
38		1-Methylnaphthalene	0.746	0.721	3.4	120	0.00
39	I	Acenaphthene-d10	1.000	1.000	0.0	123	-0.01
40		1,2,4,5-Tetrachlorobenzene	0.771	0.747	3.1	125	-0.01
41	P	Hexachlorocyclopentadiene	0.348	0.330	5.2	115	0.00
42	S	2,4,6-Tribromophenol	0.292	0.317	-8.6	145	0.00
43	C	2,4,6-Trichlorophenol	0.467	0.456	2.4	124	0.00
44		2,4,5-Trichlorophenol	0.513	0.507	1.2	126	0.00
45	S	2-Fluorobiphenyl	1.618	1.548	4.3	123	0.00
46		1,1'-Biphenyl	1.596	1.523	4.6	122	0.00
47		2-Chloronaphthalene	1.310	1.245	5.0	123	0.00

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

		Compound	AvgRF	CCRF	%Dev	Area%	Dev(min)
48		2-Nitroaniline	0.366	0.376	-2.7	133	0.00
49		Acenaphthylene	1.838	1.778	3.3	126	0.00
50		Dimethylphthalate	1.531	1.531	0.0	133	0.00
51		2,6-Dinitrotoluene	0.336	0.350	-4.2	136	0.00
52	C	Acenaphthene	1.141	1.112	2.5	127	-0.01
53		3-Nitroaniline	0.309	0.321	-3.9	142	0.00
54	P	2,4-Dinitrophenol	0.175	0.203	-16.0	158#	0.00
55		Dibenzofuran	1.858	1.833	1.3	131	0.00
56	P	4-Nitrophenol	0.253	0.279	-10.3	158#	0.00
57		2,4-Dinitrotoluene	0.441	0.491	-11.3	149	0.00
58		Fluorene	1.480	1.493	-0.9	135	0.00
59		2,3,4,6-Tetrachlorophenol	0.431	0.454	-5.3	140	0.00
60		Diethylphthalate	1.523	1.589	-4.3	143	0.00
61		4-Chlorophenyl-phenylether	0.832	0.851	-2.3	135	0.00
62		4-Nitroaniline	0.324	0.357	-10.2	159#	0.00
63		Azobenzene	1.473	1.478	-0.3	135	0.00
64	I	Phenanthrene-d10	1.000	1.000	0.0	155#	0.00
65		4,6-Dinitro-2-methylphenol	0.114	0.126	-10.5	171#	0.00
66	c	n-Nitrosodiphenylamine	0.599	0.549	8.3	141	0.00
67		4-Bromophenyl-phenylether	0.250	0.218	12.8	132	0.00
68		Hexachlorobenzene	0.292	0.277	5.1	149	0.00
69		Atrazine	0.199	0.151	24.1	123	0.00
70	C	Pentachlorophenol	0.182	0.190	-4.4	164#	0.00
71		Phenanthrene	1.053	1.013	3.8	158#	0.00
72		Anthracene	1.050	1.018	3.0	158#	0.00
73		Carbazole	0.954	0.953	0.1	173#	0.00
74		Di-n-butylphthalate	1.153	1.162	-0.8	169#	0.00
75	C	Fluoranthene	1.286	1.366	-6.2	190#	0.00
76	I	Chrysene-d12	1.000	1.000	0.0	210#	0.00
77		Benzidine	0.418	0.289	30.9#	161#	0.00
78		Pyrene	1.412	1.282	9.2	193#	0.00
79	S	Terphenyl-d14	1.216	1.151	5.3	199#	0.00
80		Butylbenzylphthalate	0.527	0.502	4.7	201#	0.00
81		Benzo(a)anthracene	1.408	1.369	2.8	214#	0.00
82		3,3'-Dichlorobenzidine	0.513	0.476	7.2	202#	0.00
83		Chrysene	1.337	1.281	4.2	211#	0.00
84		Bis(2-ethylhexyl)phthalate	0.803	0.755	6.0	198#	0.00
85	c	Di-n-octyl phthalate	1.412	1.324	6.2	196#	-0.01
86	I	Perylene-d12	1.000	1.000	0.0	201#	0.00
87		Indeno(1,2,3-cd)pyrene	1.531	1.445	5.6	197#	-0.01
88		Benzo(b)fluoranthene	1.261	1.261	0.0	207#	0.00
89		Benzo(k)fluoranthene	1.204	1.171	2.7	209#	0.00
90	C	Benzo(a)pyrene	1.122	1.106	1.4	207#	0.00
91		Dibenzo(a,h)anthracene	1.259	1.196	5.0	197#	0.00
92		Benzo(g,h,i)perylene	1.277	1.200	6.0	195#	0.00

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Compound	AvgRF	CCRF	%Dev Area%	Dev(min)

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