

4.3. ELECTRON DIFFRACTION

Table 4.3.1.2. Atomic scattering amplitudes (\AA) for electrons for ionized atoms (cont.)

Element Z Method	Ra ²⁺ 88 *DS	Ac ³⁺ 89 *DS	U ³⁺ 92 *DS	U ⁴⁺ 92 *DS	U ⁶⁺ 92 *DS
0.00					
0.01					
0.02					
0.03					
0.04	40.04	54.00	54.02	68.15	96.83
0.05	29.19	37.78	37.81	46.56	64.49
0.06	23.23	28.91	28.95	34.80	46.89
0.07	19.57	23.53	23.57	27.67	36.26
0.08	17.14	19.98	20.03	23.01	29.33
0.09	15.42	17.51	17.57	19.78	24.56
0.10	14.12	15.70	15.76	17.44	21.12
0.11	13.11	14.31	14.39	15.67	18.55
0.12	12.291	13.217	13.300	14.296	16.573
0.13	11.602	12.324	12.416	13.192	15.010
0.14	11.008	11.577	11.679	12.287	13.749
0.15	10.486	10.939	11.050	11.528	12.709
0.16	10.018	10.382	10.503	10.879	11.837
0.17	9.592	9.889	10.018	10.314	11.093
0.18	9.200	9.446	9.583	9.816	10.451
0.19	8.836	9.042	9.188	9.371	9.889
0.20	8.495	8.671	8.824	8.967	9.391
0.22	7.873	8.008	8.174	8.261	8.544
0.24	7.315	7.427	7.602	7.655	7.843
0.25	7.057	7.161	7.340	7.380	7.534
0.26	6.811	6.909	7.091	7.122	7.247
0.28	6.355	6.444	6.629	6.647	6.729
0.30	5.940	6.022	6.208	6.219	6.273
0.32	5.563	5.639	5.824	5.830	5.865
0.34	5.219	5.291	5.472	5.475	5.497
0.35	5.059	5.128	5.307	5.309	5.327
0.36	4.906	4.973	5.149	5.151	5.164
0.38	4.621	4.683	4.853	4.853	4.861
0.40	4.360	4.417	4.580	4.580	4.584
0.42	4.122	4.174	4.329	4.328	4.330
0.44	3.904	3.951	4.098	4.097	4.096
0.45	3.801	3.847	3.989	3.988	3.987
0.46	3.703	3.747	3.885	3.883	3.881
0.48	3.518	3.558	3.688	3.686	3.683
0.50	3.348	3.385	3.506	3.504	3.500
0.55	2.975	3.005	3.107	3.106	3.100
0.60	2.664	2.689	2.776	2.774	2.768
0.65	2.400	2.421	2.496	2.494	2.489
0.70	2.174	2.193	2.258	2.256	2.252
0.80	1.808	1.824	1.875	1.874	1.872
0.90	1.527	1.541	1.583	1.582	1.582
1.00	1.307	1.319	1.354	1.354	1.354
1.10	1.132	1.142	1.171	1.172	1.172
1.20	0.991	0.999	1.024	1.025	1.025
1.30	0.874	0.882	0.904	0.904	0.905
1.40	0.778	0.784	0.804	0.804	0.804
1.50	0.696	0.702	0.720	0.720	0.720
1.60	0.626	0.632	0.648	0.648	0.648
1.70	0.566	0.571	0.586	0.586	0.586
1.80	0.513	0.518	0.533	0.533	0.533
1.90	0.467	0.472	0.486	0.486	0.486
2.00	0.427	0.431	0.444	0.444	0.444

Table 4.3.2.1. Parameters useful in electron diffraction as a function of accelerating voltage, E

E (keV)	λ	$1/\lambda$	m/m_0	v/c	σ
1	0.387629	2.57979	1.00196	0.06247	0.0081126
2	0.273961	3.65016	1.00391	0.08821	0.0057448
3	0.223579	4.47270	1.00587	0.10788	0.0046975
4	0.193530	5.16715	1.00783	0.12439	0.0040741
5	0.173015	5.77986	1.00978	0.13887	0.0036493
6	0.157863	6.33460	1.01174	0.15191	0.0033361
7	0.146082	6.84548	1.01370	0.16384	0.0030931
8	0.136581	7.32168	1.01566	0.17490	0.0028975
9	0.128707	7.76958	1.01761	0.18524	0.0027358
10	0.122043	8.19383	1.01957	0.19498	0.0025991
15	0.099407	10.05963	1.02935	0.23711	0.0021374
20	0.085882	11.64383	1.03914	0.27186	0.0018641
25	0.076632	13.04940	1.04892	0.30184	0.0016790
30	0.069789	14.32899	1.05871	0.32837	0.0015433
35	0.064459	15.51381	1.06849	0.35227	0.0014386
40	0.060153	16.62414	1.07828	0.37406	0.0013548
45	0.056580	17.67403	1.08806	0.39410	0.0012859
50	0.053551	18.67366	1.09784	0.41268	0.0012280
55	0.050941	19.63072	1.10763	0.43000	0.0011786
60	0.048659	20.55115	1.11741	0.44622	0.0011357
65	0.046642	21.43968	1.12720	0.46147	0.0010982
70	0.044843	22.30012	1.13698	0.47586	0.0010650
75	0.043223	23.13560	1.14677	0.48948	0.0010354
80	0.041756	23.94874	1.15655	0.50239	0.0010087
85	0.040418	24.74173	1.16634	0.51467	0.0009847
90	0.039190	25.51646	1.17612	0.52637	0.0009628
95	0.038060	26.27454	1.18591	0.53754	0.0009428
100	0.037013	27.01738	1.19569	0.54822	0.0009244
120	0.033491	29.85866	1.23483	0.58667	0.0008638
140	0.030739	32.53222	1.27397	0.61956	0.0008180
160	0.028509	35.07642	1.31310	0.64810	0.0007820
180	0.026654	37.51759	1.35224	0.67314	0.0007529
200	0.025079	39.87466	1.39138	0.69531	0.0007289
250	0.021986	45.48412	1.48922	0.74101	0.0006839
300	0.019687	50.79517	1.58707	0.77652	0.0006526
350	0.017891	55.89295	1.68491	0.80483	0.0006297
400	0.016439	60.83109	1.78276	0.82786	0.0006122
450	0.015233	65.64563	1.88060	0.84691	0.0005984
500	0.014212	70.36195	1.97845	0.86286	0.0005873
550	0.013334	74.99858	2.07629	0.87638	0.0005783
600	0.012568	79.56945	2.17414	0.88794	0.0005707
650	0.011893	84.08529	2.27198	0.89793	0.0005644
700	0.011292	88.55452	2.36983	0.90661	0.0005590
750	0.010755	92.98385	2.46767	0.91421	0.0005543
800	0.010269	97.37874	2.56552	0.92091	0.0005503
850	0.009829	101.74364	2.66336	0.92684	0.0005468
900	0.009427	106.08226	2.76121	0.93212	0.0005437
950	0.009058	110.39769	2.85905	0.93684	0.0005410
1000	0.008719	114.69256	2.95690	0.94108	0.0005385
1100	0.008115	123.22919	3.15259	0.94836	0.0005344
1200	0.007593	131.70646	3.34828	0.95436	0.0005310
1300	0.007136	140.13516	3.54397	0.95936	0.0005282
1400	0.006733	148.52355	3.73966	0.96358	0.0005259
1500	0.006374	156.87810	3.93535	0.96718	0.0005240