

OPTICAL PROPERTIES OF SELECTED ELEMENTS

J. H. Weaver and H. P. R. Frederikse

These tables list the index of refraction n , the extinction coefficient k , and the normal incidence reflection R ($\phi = 0$) as a function of photon energy E , which is expressed in electron volts (eV). To convert the energy in eV to wavelength in μm , use $\lambda = 1.2398/E$. To compute the dielectric function $\tilde{\epsilon} = \epsilon_1 + i\epsilon_2$ from the complex index of refraction $\tilde{N} = n + ik$, use $\epsilon_1 = n^2 - k^2$ and $\epsilon_2 = 2nk$.

The optical constants in these tables are abridged from three more extensive tabulations:

- *Optical Properties of Metals* (OPM), Volumes I and II, *Physics Data*, Nr. 18-1 and 18-2, J. H. Weaver, C. Krafska, D. W. Lynch, and E. E. Koch, Fachinformationzentrum, Karlsruhe, Germany.
- *Handbook of Optical Constants* (HOC), Vol. I, 1985, and Vol. II, 1991. E. D. Palik, Ed., Academic Press, Inc.
- *American Institute of Physics Handbook* (AIPH), 3rd Edition, D. E. Gray, Ed., McGraw-Hill, New York, 1972.

The first two of these major sources provide detailed comparisons of all optical data available in the literature at the time of the compilation. For critical applications the reader should refer to the original work. References for individual metals and semiconductors are listed at the end of the tables. Generally, tabulated values

for the optical properties are accurate to better than 10%. Data in parentheses are extrapolated or interpolated values. For most elements the spectral range covered is from the far infrared (0.010 or 0.10 eV) to the far ultraviolet (10, 30 or 300 eV). The intervals between successive energies in the tables are chosen in such a way that the major spectral features are preserved.

Very small values of k are expressed in exponential notation, e.g., 1.23E-5 means 1.23×10^{-5} .

The following table is convenient for identifying the energy entries in these tables with the corresponding wavelengths:

λ	E/eV	λ	E/eV
1 mm	0.00124	6000 Å	2.066
500 μm	0.00248	5000 Å	2.480
100 μm	0.01240	4000 Å	3.100
50 μm	0.02480	3000 Å	4.133
10 μm	0.12398	2000 Å	6.199
5 μm	0.24797	1000 Å	12.398
1 μm	1.240	400 Å	30.996

Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$
<i>Aluminium¹</i>				2.200	1.018	6.846	0.9200	14.400	0.058	0.327	0.8102
0.040	98.595	203.701	0.9923	2.400	0.826	6.283	0.9228	14.600	0.067	0.273	0.7802
0.050	74.997	172.199	0.9915	2.600	0.695	5.800	0.9238	14.800	0.086	0.211	0.7202
0.060	62.852	150.799	0.9906	2.800	0.598	5.385	0.9242	15.000	0.125	0.153	0.6119
0.070	53.790	135.500	0.9899	3.000	0.523	5.024	0.9241	15.200	0.178	0.108	0.4903
0.080	45.784	123.734	0.9895	3.200	0.460	4.708	0.9243	15.400	0.234	0.184	0.3881
0.090	39.651	114.102	0.9892	3.400	0.407	4.426	0.9245	15.600	0.280	0.073	0.3182
0.100	34.464	105.600	0.9889	3.600	0.363	4.174	0.9246	15.800	0.318	0.065	0.2694
0.125	24.965	89.250	0.9884	3.800	0.326	3.946	0.9247	16.000	0.351	0.060	0.2326
0.150	18.572	76.960	0.9882	4.000	0.294	3.740	0.9248	16.200	0.380	0.055	0.2031
0.175	14.274	66.930	0.9879	4.200	0.267	3.552	0.9248	16.400	0.407	0.050	0.1789
0.200	11.733	59.370	0.9873	4.400	0.244	3.380	0.9249	16.750	0.448	0.045	0.1460
0.250	8.586	48.235	0.9858	4.600	0.223	3.222	0.9249	17.000	0.474	0.042	0.1278
0.300	6.759	40.960	0.9844	4.800	0.205	3.076	0.9249	17.250	0.498	0.040	0.1129
0.350	5.438	35.599	0.9834	5.000	0.190	2.942	0.9244	17.500	0.520	0.038	0.1005
0.400	4.454	31.485	0.9826	6.000	0.130	2.391	0.9257	17.750	0.540	0.036	0.0899
0.500	3.072	25.581	0.9817	6.500	0.110	2.173	0.9260	18.000	0.558	0.035	0.0809
0.600	2.273	21.403	0.9806	7.000	0.095	1.983	0.9262	18.500	0.591	0.032	0.0664
0.700	1.770	18.328	0.9794	7.500	0.082	1.814	0.9265	19.000	0.620	0.030	0.0554
0.800	1.444	15.955	0.9778	8.000	0.072	1.663	0.9269	19.500	0.646	0.028	0.0467
0.900	1.264	14.021	0.9749	8.500	0.063	1.527	0.9272	20.000	0.668	0.027	0.0398
1.000	1.212	12.464	0.9697	9.000	0.056	1.402	0.9277	20.500	0.689	0.025	0.0342
1.100	1.201	11.181	0.9630	9.500	0.049	1.286	0.9282	21.000	0.707	0.024	0.0296
1.200	1.260	10.010	0.9521	10.000	0.044	1.178	0.9286	21.500	0.724	0.023	0.0258
1.300	1.468	8.949	0.9318	10.500	0.040	1.076	0.9293	22.000	0.739	0.022	0.0226
1.400	2.237	8.212	0.8852	11.000	0.036	0.979	0.9298	22.500	0.753	0.021	0.0199
1.500	2.745	8.309	0.8678	11.500	0.033	0.883	0.9283	23.000	0.766	0.021	0.0177
1.600	2.625	8.597	0.8794	12.000	0.033	0.791	0.9224	23.500	0.778	0.020	0.0157
1.700	2.143	8.573	0.8972	12.500	0.034	0.700	0.9118	24.000	0.789	0.019	0.0140
1.800	1.741	8.205	0.9069	13.000	0.038	0.609	0.8960	24.500	0.799	0.018	0.0126
1.900	1.488	7.821	0.9116	13.500	0.041	0.517	0.8789	25.000	0.809	0.018	0.0113
2.000	1.304	7.479	0.9148	14.000	0.048	0.417	0.8486	25.500	0.817	0.017	0.0102
				14.200	0.053	0.373	0.8312	26.000	0.826	0.016	0.0092

Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$
<i>Chromium⁴</i>				11.00	1.05	0.69	0.103	2.20	2.07	3.70	0.642
0.06	21.19	42.00	0.962	11.50	1.09	0.69	0.100	2.30	2.01	3.59	0.634
0.10	11.81	29.76	0.955	12.00	1.13	0.70	0.101	2.40	1.95	3.49	0.627
0.14	15.31	26.36	0.936	12.50	1.15	0.73	0.108	2.50	1.88	3.40	0.622
0.18	8.73	25.37	0.53	13.00	1.15	0.77	0.119	2.60	1.81	3.32	0.618
0.22	5.30	20.62	0.954	13.50	1.12	0.80	0.128	2.70	1.73	3.24	0.615
0.26	3.91	17.12	0.951	14.00	1.09	0.82	0.135	2.80	1.66	3.13	0.607
0.30	3.15	14.28	0.943	14.50	1.03	0.82	0.142	2.90	1.61	3.05	0.600
0.42	3.47	8.97	0.862	15.00	1.00	0.82	0.143	3.00	1.55	2.96	0.594
0.54	3.92	7.06	0.788	15.50	0.96	0.80	0.141	3.20	1.46	2.80	0.579
0.66	3.96	5.95	0.736	16.00	0.92	0.77	0.139	3.40	1.38	2.64	0.563
0.78	4.13	5.03	0.680	16.50	0.31	0.75	0.134	3.60	1.31	2.48	0.544
0.90	4.43	4.60	0.650	17.00	0.90	0.73	0.132	3.80	1.28	2.33	0.519
1.00	4.47	4.43	0.639	17.50	0.88	0.72	0.130	4.00	1.26	2.20	0.495
1.12	4.53	4.31	0.631	18.00	0.87	0.70	0.129	4.20	1.25	2.10	0.471
1.24	4.50	4.28	0.629	18.50	0.84	0.69	0.130	4.40	1.24	2.01	0.452
1.36	4.42	4.30	0.631	19.00	0.82	0.68	0.131	4.60	1.24	1.94	0.435
1.46	4.31	4.32	0.632	20.00	0.77	0.64	0.130	4.80	1.23	1.88	0.423
1.77	3.84	4.37	0.639	20.5	0.76	0.63	0.129	5.00	1.22	1.83	0.411
2.00	3.48	4.36	0.644	21.0	0.74	0.58	0.121	5.20	1.21	1.79	0.403
2.20	3.18	4.41	0.656	21.5	0.72	0.55	0.116	5.40	1.19	1.77	0.399
2.40	2.75	4.46	0.677	22.0	0.71	0.52	0.112	5.60	1.16	1.75	0.400
2.60	2.22	4.36	0.698	22.5	0.70	0.50	0.109	5.80	1.10	1.73	0.406
2.80	1.80	4.06	0.703	23.0	0.69	0.48	0.105	6.00	1.03	1.68	0.407
3.00	1.54	3.71	0.695	23.5	0.68	0.45	0.101	6.20	0.97	1.62	0.401
3.20	1.44	3.40	0.670	24.0	0.68	0.43	0.096	6.40	0.94	1.53	0.386
3.40	1.39	3.24	0.657	24.5	0.67	0.39	0.089	6.60	0.91	1.46	0.368
3.60	1.26	3.12	0.661	25.0	0.68	0.36	0.080	6.80	0.91	1.38	0.345
3.80	1.12	2.95	0.660	25.5	0.68	0.33	0.072	7.00	0.91	1.32	0.326
4.00	1.02	2.76	0.651	26.0	0.70	0.31	0.063	7.00	0.91	1.26	0.305
4.20	0.94	2.58	0.639	26.5	0.71	0.28	0.055	7.40	0.92	1.21	0.286
4.40	0.90	2.42	0.620	27.0	0.72	0.26	0.048	7.60	0.93	1.17	0.269
4.50	0.89	2.35	0.607	27.5	0.73	0.25	0.043	7.80	0.94	1.13	0.253
4.60	0.88	2.28	0.598	28.0	0.75	0.23	0.037	8.00	0.95	1.09	0.239
4.70	0.86	2.21	0.586	29.0	0.77	0.22	0.032	<i>Cobalt, single crystal, $\vec{E} \perp c^s$</i>			
4.80	0.86	2.13	0.572	30.0	0.78	0.21	0.030	0.10	5.83	32.36	0.979
4.90	0.86	2.07	0.557	<i>Cobalt, single crystal, $\vec{E} \parallel c^s$</i>				0.15	4.24	21.37	0.965
5.00	0.85	2.01	0.542	0.10	6.71	37.87	0.982	0.20	3.87	15.53	0.042
5.10	0.86	1.94	0.523	0.15	4.66	25.47	0.973	0.30	4.34	10.01	0.865
5.20	0.87	1.87	0.503	0.20	3.55	18.78	0.962	0.40	4.66	7.39	0.785
5.40	0.93	1.80	0.466	0.25	3.98	14.59	0.933	0.50	5.17	5.75	0.709
5.60	0.95	1.74	0.443	0.30	4.04	12.16	0.907	0.60	5.77	5.17	0.682
5.80	0.97	1.74	0.437	0.40	4.24	9.13	0.847	0.70	6.15	5.20	0.685
6.00	0.94	1.73	0.444	0.50	4.41	7.19	0.782	0.80	6.08	5.61	0.702
6.20	0.89	1.69	0.446	0.60	4.91	6.13	0.729	0.90	5.57	5.93	0.715
6.40	0.85	1.66	0.447	0.70	5.24	5.85	0.713	1.00	4.83	5.94	0.721
6.60	0.80	1.59	0.444	0.80	5.17	5.89	0.716	1.10	4.31	5.60	0.711
6.80	0.75	1.51	0.439	0.90	4.94	5.95	0.720	1.20	4.02	5.34	0.701
7.00	0.74	1.45	0.425	1.00	4.46	5.86	0.722	1.30	3.78	5.16	0.694
7.20	0.71	1.39	0.414	1.10	4.07	5.61	0.715	1.40	3.55	5.05	0.692
7.40	0.69	1.33	0.404	1.20	3.81	5.36	0.706	1.50	3.26	4.93	0.692
7.60	0.66	1.23	0.378	1.30	3.60	5.20	0.701	1.60	3.03	4.74	0.687
7.80	0.67	1.15	0.347	1.40	3.37	5.09	0.701	1.70	2.83	4.60	0.684
8.00	0.68	1.07	0.315	1.50	3.10	4.96	0.701	1.80	2.61	4.45	0.683
8.20	0.71	1.00	0.278	1.60	2.84	4.77	0.697	1.90	2.41	4.27	0.677
8.50	0.74	0.92	0.235	1.70	2.66	4.57	0.690	2.00	2.25	4.09	0.670
9.0	0.83	0.81	0.170	1.80	2.45	4.41	0.687	2.10	2.13	3.89	0.659
9.50	0.92	0.74	0.132	1.90	2.31	4.18	0.675	2.20	2.04	3.72	0.646
10.00	0.98	0.73	0.120	2.00	2.21	4.00	0.664	2.30	1.99	3.56	0.632
10.50	1.01	0.72	0.112	2.10	2.13	3.85	0.654	2.40	1.95	3.44	0.620

Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$	
2.50	1.90	3.34	0.611	5.60	1.18	1.74	0.391	58.00	0.96	0.11	0.004	
2.60	1.86	3.26	0.605	5.80	1.10	1.67	0.389	59.00	0.97	0.11	0.003	
2.70	1.79	3.19	0.602	6.00	1.04	1.59	0.380	60.00	0.97	0.11	0.003	
2.80	1.72	3.11	0.596	6.50	0.96	1.37	0.329	61.00	0.97	0.11	0.003	
2.90	1.66	3.03	0.591	7.00	0.97	1.20	0.271	62.00	0.97	0.11	0.003	
3.00	1.60	2.94	0.586	7.50	1.00	1.09	0.230	63.00	0.96	0.10	0.003	
3.20	1.50	2.78	0.571	8.00	1.03	1.03	0.206	64.00	0.96	0.10	0.003	
3.40	1.42	2.62	0.553	8.50	1.03	0.98	0.189	65.00	0.97	0.10	0.003	
3.60	1.36	2.47	0.533	9.00	1.03	0.92	0.171	66.00	0.97	0.10	0.003	
3.80	1.33	2.33	0.511	9.50	1.03	0.87	0.154	67.00	0.97	0.09	0.003	
4.00	1.31	2.21	0.488	10.00	1.04	0.82	0.139	68.00	0.97	0.09	0.002	
4.20	1.28	2.12	0.471	11.00	1.07	0.75	0.118	69.00	0.97	0.09	0.002	
4.40	1.27	2.03	0.452	12.00	1.09	0.73	0.111	70.00	0.97	0.09	0.002	
4.60	1.26	1.95	0.435	13.00	1.08	0.72	0.109	75.00	0.98	0.09	0.002	
4.80	1.25	1.90	0.423	14.00	1.06	0.72	0.111	80.00	0.98	0.09	0.002	
5.00	1.24	1.84	0.411	14.50	1.03	0.72	0.111	85.00	0.97	0.09	0.002	
5.20	1.22	1.80	0.403	15.00	1.01	0.71	0.111	90.00	0.96	0.08	0.002	
5.40	1.21	1.78	0.399	15.50	0.98	0.69	0.109	<i>Gallium (liquid)⁷</i>				
5.60	1.17	1.76	0.400	16.00	0.95	0.67	0.106	1.425	2.40	9.20	0.900	
5.80	1.11	1.74	0.406	17.00	0.91	0.62	0.097	1.550	2.09	8.50	0.898	
6.00	1.04	1.69	0.407	18.00	0.89	0.56	0.084	1.771	1.65	7.60	0.898	
6.20	0.98	1.62	0.401	19.00	0.88	0.51	0.071	2.066	1.25	6.60	0.897	
6.40	0.94	1.54	0.386	20.00	0.88	0.45	0.059	2.480	0.89	5.60	0.898	
6.60	0.92	1.46	0.368	21.00	0.90	0.41	0.048	3.100	0.59	4.50	0.896	
6.80	0.91	1.38	0.345	22.00	0.92	0.38	0.040	<i>Germanium, single crystal⁸</i>				
7.00	0.91	1.32	0.326	23.00	0.94	0.37	0.035	0.01240	(4.0065)	3.00E-03	0.361	
7.20	0.91	1.26	0.305	24.00	0.96	0.37	0.035	0.01364	4.0063	2.40E-03	0.361	
7.40	0.92	1.21	0.285	25.00	0.96	0.40	0.040	0.01488	(4.0060)	1.70E-03	0.361	
7.60	0.93	1.17	0.269	26.00	0.92	0.40	0.044	0.01612	(4.0060)	1.55E-03	0.361	
7.80	0.94	1.13	0.253	27.00	0.88	0.38	0.043	0.01736	(4.0060)	1.50E-03	0.361	
<i>Copper⁶</i>				28.00	0.86	0.35	0.039	0.01860		1.50E-03		
0.10	29.69	71.57	0.980	29.00	0.85	0.30	0.032	0.01984		1.60E-03		
0.50	1.71	17.63	0.979	30.00	0.86	0.26	0.025	0.02108		1.60E-03		
1.00	0.44	8.48	0.976	31.00	0.88	0.24	0.020	0.02232		1.55E-03		
1.50	0.26	5.26	0.965	32.00	0.89	0.22	0.017	0.02356		1.53E-03		
1.70	0.22	4.43	0.958	33.00	0.90	0.21	0.015	0.02480		1.50E-03		
1.75	0.21	4.25	0.956	34.00	0.91	0.20	0.014	0.02604		1.25E-03		
1.80	0.21	4.04	0.952	35.00	0.92	0.20	0.013	0.02728		8.50E-04		
1.85	0.22	3.85	0.947	36.00	0.92	0.19	0.012	0.02852		6.50E-04		
1.90	0.21	3.67	0.943	37.00	0.92	0.19	0.011	0.02976		7.00E-04		
2.00	0.27	3.24	0.910	38.00	0.93	0.18	0.010	0.03100	3.9827	8.50E-04	0.358	
2.10	0.47	2.81	0.814	39.00	0.93	0.17	0.009	0.03224		1.55E-03		
2.20	0.83	2.60	0.673	40.00	0.93	0.17	0.009	0.03348		2.75E-03		
2.30	1.04	2.59	0.618	41.00	0.94	0.16	0.008	0.03472		3.55E-03		
2.40	1.12	2.60	0.602	42.00	0.94	0.16	0.007	0.03596	(3.9900)	3.05E-03	0.359	
2.60	1.15	2.50	0.577	43.00	0.94	0.15	0.007	0.03720		2.75E-03		
2.80	1.17	2.36	0.545	44.00	0.95	0.15	0.007	0.03844		2.70E-03		
3.00	1.18	2.21	0.509	45.00	0.95	0.15	0.006	0.03968	(3.9930)	2.90E-03	0.359	
3.20	1.23	2.07	0.468	46.00	0.95	0.15	0.006	0.04092		2.95E-03		
3.40	1.27	1.95	0.434	47.00	0.95	0.14	0.006	0.04215		3.20E-03		
3.60	1.31	1.87	0.407	48.00	0.95	0.14	0.006	0.04339		6.30E-03		
3.80	1.34	1.81	0.387	49.00	0.95	0.14	0.005	0.04463		3.40E-03		
4.00	1.34	1.72	0.364	50.00	0.95	0.13	0.005	0.04587	(3.9955)	2.50E-03	0.360	
4.20	1.42	1.64	0.336	51.00	0.95	0.13	0.005	0.04711		2.10E-03		
4.40	1.49	1.64	0.329	52.00	0.95	0.13	0.005	0.04835		2.00E-03		
4.60	1.52	1.67	0.334	53.00	0.96	0.12	0.004	0.04959		8.00E-04		
4.80	1.53	1.71	0.345	54.00	0.96	0.12	0.004	0.05083		1.40E-03		
5.00	1.47	1.78	0.366	55.00	0.96	0.12	0.004	0.05207		1.35E-03		
5.20	1.38	1.80	0.380	56.00	0.96	0.11	0.004	0.05331		1.10E-03		
5.40	1.28	1.78	0.389	57.00	0.96	0.11	0.004	0.05455		8.00E-04		

Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)	Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)	Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)
19.50	0.75	0.47	0.086	6.15	0.376	0.522	0.306	4.24	1.85	2.14	0.417
19.67	0.75	0.47	0.085	6.45	0.408	0.460	0.256	4.36	1.85	2.08	0.406
19.83	0.75	0.46	0.084	6.75	0.440	0.407	0.214	4.49	1.86	2.03	0.395
20.00	0.74	0.45	0.083	7.05	0.466	0.364	0.183	4.61	1.85	1.99	0.388
20.17	0.74	0.44	0.081	7.35	0.492	0.320	0.155	4.74	1.84	1.94	0.378
20.33	0.74	0.44	0.081	7.65	0.517	0.282	0.131	4.86	1.83	1.91	0.372
20.50	0.74	0.42	0.080	7.95	0.545	0.246	0.109	4.98	1.82	1.86	0.362
20.67	0.73	0.43	0.079	8.25	0.572	0.214	0.091	5.11	1.82	1.82	0.354
20.83	0.73	0.42	0.078	8.55	0.601	0.189	0.075	5.23	1.81	1.79	0.348
21.00	0.73	0.41	0.077	8.85	0.624	0.163	0.063	5.36	1.78	1.76	0.342
21.17	0.72	0.40	0.076	9.15	0.657	0.144	0.050	5.48	1.74	1.73	0.337
21.33	0.72	0.39	0.074	9.45	0.680	0.130	0.042	5.60	1.73	1.70	0.331
21.50	0.72	0.38	0.073	9.75	0.708	0.119	0.034	5.73	1.72	1.67	0.325
21.67	0.72	0.38	0.071	10.1	0.726	0.108	0.029	5.85	1.70	1.64	0.319
21.83	0.72	0.37	0.070	10.4	0.743	0.102	0.025	5.98	1.67	1.61	0.313
22.00	0.72	0.36	0.068	10.6	0.753	0.080	0.022	6.10	1.63	1.58	0.307
22.17	0.71	0.35	0.067	<i>Magnesium (evaporated)¹³</i>				6.22	1.62	1.55	0.301
22.33	0.72	0.34	0.064	2.145	0.48	3.71	0.880	6.35	1.59	1.52	0.295
22.50	0.72	0.34	0.063	2.270	0.57	3.47	0.843	6.47	1.55	1.50	0.292
22.67	0.72	0.33	0.062	2.522	0.53	2.92	0.805	6.60	1.48	1.47	0.288
22.83	0.72	0.32	0.059	2.845	0.52	2.65	0.777	<i>Mercury (liquid)¹⁵</i>			
23.00	0.72	0.31	0.058	3.064	0.52	2.05	0.681	0.2	13.99	14.27	0.869
23.17	0.72	0.30	0.056	5.167	0.10	1.60	0.894	0.3	11.37	11.95	0.846
23.33	0.72	0.29	0.054	5.636	0.15	1.50	0.832	0.4	9.741	10.65	0.830
23.50	0.73	0.28	0.050	6.200	0.20	1.40	0.765	0.5	8.528	9.805	0.818
23.67	0.73	0.28	0.049	6.889	0.25	1.30	0.693	0.6	7.574	9.195	0.808
23.83	0.74	0.27	0.047	7.750	0.20	1.20	0.722	0.8	6.086	8.312	0.796
24.00	0.74	0.27	0.045	8.857	0.15	0.95	0.730	1.0	4.962	7.643	0.789
24.17	0.74	0.26	0.044	10.335	0.25	0.40	0.419	1.2	4.050	7.082	0.786
24.33	0.74	0.26	0.043	<i>Manganese¹⁴</i>				1.4	3.324	6.558	0.785
24.50	0.74	0.25	0.042	0.64	3.89	5.95	0.738	1.6	2.746	6.054	0.783
24.67	0.75	0.25	0.040	0.77	3.78	5.41	0.710	1.8	2.284	5.582	0.782
24.83	0.75	0.24	0.039	0.89	3.65	5.02	0.688	2.0	1.910	5.150	0.782
25.00	0.75	0.24	0.038	1.02	3.48	4.74	0.673	2.2	1.620	4.751	0.780
26.00	0.76	0.21	0.031	1.14	3.30	4.53	0.662	2.4	1.384	4.407	0.779
27.00	0.78	0.18	0.026	1.26	3.10	4.35	0.653	2.6	1.186	4.090	0.779
28.00	0.79	0.16	0.021	1.39	2.97	4.18	0.643	2.8	1.027	3.802	0.779
29.00	0.81	0.14	0.017	1.51	2.83	4.03	0.634	3.0	0.898	3.538	0.777
30.00	0.82	0.13	0.014	1.64	2.70	3.91	0.627	3.2	0.798	3.294	0.773
<i>Lithium¹²</i>				1.76	2.62	3.78	0.617	3.4	0.713	3.074	0.770
0.14	0.659	38.0	0.998	1.88	2.56	3.65	0.606	3.6	0.644	2.860	0.763
0.54	0.661	12.6	0.984	2.01	2.51	3.54	0.596	3.8	0.589	2.665	0.755
0.75	0.561	7.68	0.963	2.13	2.47	3.43	0.585	4.0	0.542	2.502	0.749
1.05	0.448	5.58	0.946	2.26	2.39	3.33	0.577	4.2	0.507	2.341	0.738
1.35	0.338	4.36	0.935	2.38	2.32	3.23	0.567	4.4	0.477	2.195	0.727
1.65	0.265	3.55	0.925	2.50	2.25	3.14	0.559	4.6	0.452	2.058	0.715
1.95	0.221	2.94	0.913	2.63	2.19	3.06	0.552	4.8	0.431	1.929	0.701
2.25	0.206	2.48	0.892	2.75	2.11	2.98	0.545	5.0	0.414	1.806	0.685
2.55	0.217	2.11	0.854	2.88	2.06	2.90	0.536	5.2	0.401	1.687	0.666
2.85	0.247	1.82	0.797	3.00	2.00	2.82	0.528	5.4	0.394	1.569	0.642
3.15	0.304	1.60	0.715	3.12	1.96	2.74	0.518	5.6	0.386	1.454	0.617
3.45	0.334	1.45	0.656	3.25	1.92	2.67	0.509	5.7	0.386	1.396	0.601
3.75	0.345	1.32	0.611	3.37	1.89	2.59	0.498	5.8	0.386	1.341	0.585
4.05	0.346	1.21	0.578	3.50	1.89	2.51	0.484	5.9	0.385	1.287	0.569
4.35	0.333	1.11	0.557	3.62	1.87	2.45	0.475	6.0	0.386	1.232	0.551
4.65	0.317	1.01	0.540	3.74	1.86	2.38	0.463	6.1	0.388	1.176	0.531
4.95	0.302	0.906	0.520	3.87	1.86	2.32	0.451	6.2	0.390	1.118	0.510
5.25	0.299	0.795	0.484	3.99	1.86	2.25	0.438	6.3	0.399	1.058	0.481
5.55	0.310	0.688	0.434	4.12	1.86	2.19	0.427	6.4	0.412	1.002	0.450
5.85	0.342	0.594	0.365					6.5	0.428	0.949	0.418

Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$
36.00	0.82	0.34	0.043	6.60	1.01	1.40	0.325	25.00	0.89	0.42	0.050
37.00	0.81	0.30	0.038	6.80	1.02	1.35	0.308	26.00	0.88	0.39	0.046
38.00	0.81	0.27	0.033	7.00	1.03	1.30	0.291	27.00	0.87	0.37	0.042
39.00	0.82	0.25	0.029	7.20	1.03	1.27	0.282	28.00	0.87	0.35	0.040
40.00	0.83	0.23	0.025	7.40	1.03	1.24	0.273	29.00	0.86	0.34	0.037
<i>Nickel</i> ¹⁷				7.60	1.02	1.22	0.265	30.00	0.86	0.32	0.034
0.10	9.54	45.82	0.983	7.80	1.01	1.18	0.256	35.00	0.86	0.24	0.022
0.15	5.45	30.56	0.978	8.00	1.01	1.15	0.248	40.00	0.87	0.18	0.014
0.20	4.12	22.48	0.969	8.20	1.00	1.13	0.242	45.00	0.88	0.13	0.008
0.25	4.25	17.68	0.950	8.40	0.99	1.11	0.235	50.00	0.92	0.10	0.004
0.30	4.19	15.05	0.934	8.60	0.98	1.08	0.228	60.00	0.96	0.08	0.002
0.35	4.03	13.05	0.918	8.80	0.97	1.05	0.220	65.00	0.98	0.09	0.002
0.40	3.84	11.43	0.900	9.00	0.97	1.01	0.211	68.00	0.96	0.12	0.004
0.50	4.03	9.64	0.864	9.20	0.96	0.99	0.203	70.00	0.94	0.11	0.004
0.60	3.84	8.35	0.835	9.40	0.95	0.96	0.194	75.00	0.94	0.09	0.003
0.70	3.59	7.48	0.813	9.60	0.95	0.93	0.185	80.00	0.94	0.07	0.002
0.80	3.38	6.82	0.794	9.80	0.95	0.89	0.175	90.00	0.94	0.06	0.002
0.90	3.18	6.23	0.774	10.00	0.95	0.87	0.166	<i>Niobium</i> ¹⁸			
1.00	3.06	5.74	0.753	10.20	0.95	0.83	0.155	0.12	15.99	53.20	0.979
1.10	2.97	5.38	0.734	10.40	0.95	0.80	0.145	0.20	7.25	34.14	0.976
1.20	2.85	5.10	0.721	10.60	0.97	0.76	0.129	0.24	5.47	28.88	0.975
1.30	2.74	4.85	0.708	10.80	0.99	0.75	0.123	0.28	4.26	24.95	0.974
1.40	2.65	4.63	0.695	11.00	1.01	0.73	0.115	0.35	3.11	20.03	0.970
1.50	2.53	4.47	0.688	11.25	1.04	0.72	0.111	0.45	2.28	15.58	0.964
1.60	2.43	4.31	0.679	11.50	1.05	0.71	0.109	0.55	1.83	12.67	0.956
1.70	2.28	4.18	0.677	11.75	1.07	0.71	0.108	0.65	1.57	10.59	0.947
1.80	2.14	4.01	0.670	12.00	1.07	0.71	0.108	0.75	1.41	9.00	0.935
1.90	2.02	3.82	0.659	12.25	1.07	0.71	0.107	0.85	1.35	7.74	0.918
2.00	1.92	3.65	0.649	12.50	1.08	0.71	0.106	0.95	1.35	6.70	0.893
2.10	1.85	3.48	0.634	12.75	1.08	0.71	0.106	1.05	1.44	5.86	0.857
2.20	1.80	3.33	0.620	13.00	1.08	0.71	0.105	1.15	1.55	5.18	0.814
2.30	1.75	3.19	0.605	13.25	1.08	0.71	0.105	1.25	1.65	4.63	0.768
2.40	1.71	3.06	0.590	13.50	1.07	0.70	0.105	1.35	1.76	4.13	0.715
2.50	1.67	2.93	0.575	13.75	1.07	0.70	0.105	1.45	1.95	3.68	0.650
2.60	1.65	2.81	0.557	14.00	1.07	0.71	0.106	1.55	2.15	3.37	0.595
2.70	1.64	2.71	0.542	14.25	1.06	0.70	0.106	1.65	2.36	3.13	0.552
2.80	1.63	2.61	0.525	14.50	1.05	0.70	0.106	1.75	2.54	2.99	0.527
2.90	1.62	2.52	0.509	14.75	1.04	0.70	0.107	1.85	2.69	2.89	0.510
3.00	1.61	2.44	0.495	15.00	1.03	0.70	0.107	1.95	2.82	2.86	0.505
3.10	1.61	2.36	0.480	15.25	1.02	0.69	0.106	2.05	2.89	2.87	0.505
3.20	1.61	2.30	0.467	15.50	1.01	0.69	0.105	2.15	2.92	2.87	0.505
3.30	1.61	2.23	0.454	15.75	1.00	0.68	0.104	2.25	2.93	2.87	0.505
3.40	1.62	2.17	0.441	16.00	0.99	0.67	0.103	2.35	2.92	2.88	0.506
3.50	1.63	2.11	0.428	16.50	0.98	0.66	0.101	2.45	2.89	2.90	0.509
3.60	1.64	2.07	0.416	17.00	0.96	0.64	0.098	2.55	2.83	2.92	0.512
3.70	1.66	2.02	0.405	17.50	0.94	0.63	0.096	2.65	2.74	2.90	0.511
3.80	1.69	1.99	0.397	18.00	0.92	0.61	0.092	2.75	2.66	2.86	0.507
3.90	1.72	1.98	0.393	18.50	0.91	0.58	0.087	2.85	2.58	2.80	0.500
4.00	1.73	1.98	0.392	19.00	0.90	0.56	0.082	3.00	2.51	2.68	0.485
4.20	1.74	2.01	0.396	19.50	0.90	0.54	0.077	3.10	2.48	2.60	0.475
4.40	1.71	2.06	0.409	20.00	0.89	0.51	0.071	3.20	2.45	2.53	0.465
4.60	1.63	2.09	0.421	20.50	0.89	0.49	0.066	3.30	2.44	2.45	0.453
4.80	1.53	2.11	0.435	21.00	0.90	0.47	0.061	3.40	2.46	2.38	0.442
5.00	1.40	2.10	0.449	21.50	0.91	0.46	0.057	3.50	2.48	2.33	0.435
5.20	1.27	2.04	0.454	22.00	0.91	0.45	0.055	3.60	2.52	2.29	0.428
5.40	1.16	1.94	0.449	22.50	0.91	0.44	0.053	3.70	2.56	2.27	0.426
5.60	1.09	1.83	0.435	23.00	0.92	0.44	0.051	3.80	2.59	2.28	0.427
5.80	1.04	1.73	0.417	23.50	0.91	0.44	0.052	3.90	2.62	2.29	0.429
6.20	1.00	1.54	0.371	24.00	0.90	0.43	0.051	4.00	2.64	2.33	0.434
6.40	1.01	1.46	0.345	24.50	0.90	0.43	0.051	4.20	2.64	2.42	0.447

Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)	Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)	Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)
4.40	2.53	2.56	0.467	20.60	0.71	0.55	0.119	2.20	4.58	1.62	0.457
4.60	2.39	2.56	0.470	21.00	0.72	0.50	0.100	2.30	4.84	1.76	0.479
4.80	2.32	2.52	0.465	21.60	0.75	0.43	0.075	2.40	5.10	2.01	0.506
5.00	2.26	2.57	0.475	22.00	0.78	0.40	0.063	2.50	5.28	2.38	0.532
5.20	2.16	2.62	0.487	22.60	0.82	0.35	0.045	2.60	5.36	2.82	0.557
5.40	2.00	2.68	0.505	23.00	0.85	0.33	0.038	2.70	5.30	3.29	0.580
5.60	1.81	2.67	0.518	23.60	0.88	0.30	0.029	2.80	5.07	3.78	0.603
5.80	1.63	2.60	0.522	24.00	0.91	0.29	0.025	2.90	4.65	4.18	0.624
6.00	1.49	2.49	0.520	24.60	0.94	0.28	0.022	3.00	4.05	4.40	0.639
6.20	1.38	2.38	0.512	25.00	0.96	0.27	0.020	3.20	3.29	3.96	0.614
6.40	1.31	2.25	0.496	25.60	0.99	0.26	0.018	3.40	2.93	3.79	0.607
6.60	1.26	2.14	0.480	26.00	1.00	0.26	0.017	3.60	2.75	3.45	0.577
6.80	1.24	2.04	0.460	26.60	1.03	0.25	0.016	3.80	2.73	3.32	0.562
7.00	1.23	1.96	0.441	27.00	1.04	0.25	0.015	4.00	2.71	3.34	0.565
7.20	1.22	1.91	0.430	27.60	1.06	0.25	0.015	4.20	2.53	3.44	0.584
7.40	1.20	1.88	0.427	28.00	1.08	0.24	0.015	4.40	2.24	3.44	0.599
7.60	1.14	1.85	0.430	28.60	1.11	0.24	0.016	4.60	2.01	3.31	0.598
7.80	1.07	1.78	0.428	29.00	1.13	0.25	0.017	4.80	1.88	3.19	0.592
8.00	1.02	1.69	0.412	29.60	1.16	0.26	0.020	5.00	1.74	3.12	0.596
8.20	1.00	1.60	0.390	30.00	1.18	0.28	0.023	5.20	1.58	3.00	0.597
8.40	0.99	1.51	0.365	31.00	1.18	0.31	0.026	5.40	1.46	2.88	0.593
8.60	0.99	1.43	0.340	32.00	1.20	0.34	0.031	5.60	1.36	2.77	0.589
8.70	0.99	1.39	0.328	33.00	1.21	0.38	0.038	5.80	1.27	2.65	0.582
8.80	1.00	1.36	0.315	34.00	1.20	0.42	0.044	6.00	1.20	2.54	0.575
9.00	1.01	1.29	0.290	35.20	1.17	0.47	0.051	6.20	1.13	2.44	0.571
9.20	1.04	1.22	0.265	36.00	1.15	0.50	0.056	6.40	1.06	2.33	0.562
9.40	1.07	1.18	0.245	37.50	1.07	0.53	0.064	6.60	1.01	2.21	0.548
9.60	1.10	1.13	0.227	39.50	0.95	0.50	0.063	6.80	0.97	2.11	0.532
9.80	1.13	1.09	0.209	40.50	0.92	0.47	0.059	7.00	0.95	2.00	0.514
10.00	1.18	1.05	0.194					7.20	0.92	1.91	0.497
10.20	1.23	1.04	0.187	<i>Osmium (Polycrystalline)⁹</i>				7.40	0.91	1.81	0.476
10.40	1.27	1.04	0.185	0.10	4.08	50.23	0.994	7.60	0.90	1.72	0.451
10.60	1.30	1.06	0.190	0.15	2.90	33.60	0.990	7.80	0.90	1.63	0.426
10.80	1.32	1.08	0.195	0.20	2.44	25.11	0.985	8.00	0.91	1.55	0.400
11.00	1.32	1.10	0.200	0.25	2.35	19.99	0.977	8.20	0.91	1.48	0.375
11.20	1.31	1.12	0.204	0.30	2.23	16.54	0.969	8.40	0.94	1.40	0.344
11.40	1.30	1.13	0.207	0.35	2.33	14.06	0.955	8.60	0.96	1.34	0.319
11.60	1.28	1.13	0.209	0.40	2.45	12.32	0.940	8.80	0.98	1.29	0.296
11.80	1.27	1.13	0.210	0.45	2.43	11.02	0.927	9.00	1.01	1.24	0.274
12.00	1.25	1.12	0.209	0.50	2.41	9.97	0.913	9.20	1.04	1.19	0.255
12.40	1.24	1.10	0.204	0.55	2.33	9.12	0.901	9.40	1.08	1.16	0.238
12.80	1.24	1.09	0.200	0.60	2.21	8.37	0.890	9.60	1.10	1.14	0.229
13.20	1.24	1.09	0.201	0.65	2.11	7.68	0.877	9.80	1.13	1.11	0.217
13.60	1.23	1.12	0.208	0.70	2.02	7.04	0.862	10.00	1.16	1.10	0.209
14.00	1.20	1.13	0.216	0.75	2.00	6.46	0.842	10.20	1.19	1.08	0.203
14.40	1.16	1.15	0.225	0.80	2.00	5.95	0.820	10.30	1.20	1.08	0.201
14.80	1.11	1.16	0.234	0.85	2.01	5.51	0.796	10.40	1.22	1.08	0.200
15.00	1.08	1.16	0.238	0.90	2.03	5.10	0.769	10.50	1.23	1.09	0.201
15.60	0.99	1.14	0.247	0.95	2.05	4.74	0.742	10.60	1.24	1.10	0.203
16.00	0.92	1.11	0.250	1.00	2.09	4.41	0.712	10.80	1.25	1.11	0.206
16.60	0.85	1.04	0.245	1.10	2.15	3.84	0.651	11.00	1.24	1.13	0.213
17.00	0.80	0.99	0.240	1.20	2.16	3.35	0.592	11.20	1.23	1.14	0.217
17.20	0.79	0.96	0.236	1.30	2.25	2.77	0.506	11.40	1.19	1.15	0.223
17.40	0.77	0.93	0.230	1.40	2.49	2.23	0.419	11.60	1.17	1.12	0.216
17.80	0.75	0.87	0.217	1.50	2.84	1.80	0.369	11.80	1.16	1.10	0.211
18.00	0.74	0.85	0.209	1.60	3.36	1.62	0.379	12.00	1.15	1.08	0.205
18.60	0.73	0.77	0.185	1.70	3.70	1.75	0.411	12.40	1.14	1.03	0.191
19.00	0.72	0.72	0.170	1.80	3.78	1.83	0.423	12.80	1.15	1.01	0.183
19.60	0.72	0.66	0.150	1.90	3.81	1.75	0.418	13.20	1.16	0.98	0.174
20.00	0.72	0.62	0.137	2.00	3.98	1.60	0.418	13.60	1.17	0.97	0.170
				2.10	4.26	1.54	0.432				

Energy (eV)	n	k	$R(\phi=0)$	Energy (eV)	n	k	$R(\phi=0)$	Energy (eV)	n	k	$R(\phi=0)$
14.00	1.17	0.96	0.169	0.56	3.92	10.49	0.883	9.50	1.12	0.65	0.089
14.40	1.16	0.94	0.165	0.60	3.80	9.96	0.876	10.00	1.14	0.65	0.088
14.80	1.16	0.91	0.156	0.72	3.51	8.70	0.854	10.50	1.16	0.65	0.087
15.20	1.17	0.89	0.148	0.80	3.35	8.06	0.840	11.00	1.18	0.64	0.086
15.60	1.20	0.86	0.140	1.00	2.99	6.89	0.811	11.50	1.19	0.65	0.087
16.00	1.25	0.87	0.140	1.10	2.81	6.46	0.800	12.00	1.20	0.66	0.089
16.40	1.28	0.90	0.147	1.20	2.65	6.10	0.790	12.50	1.19	0.67	0.091
16.80	1.28	0.94	0.157	1.30	2.50	5.78	0.781	13.00	1.18	0.67	0.091
17.20	1.27	0.97	0.167	1.40	2.34	5.50	0.774	13.50	1.18	0.67	0.092
17.60	1.26	1.01	0.178	1.50	2.17	5.22	0.767	14.00	1.17	0.67	0.093
18.00	1.23	1.04	0.189	1.60	2.08	4.95	0.755	14.50	1.15	0.68	0.095
18.40	1.19	1.08	0.200	1.70	2.00	4.72	0.745	15.00	1.13	0.69	0.098
18.80	1.14	1.10	0.210	1.80	1.92	4.54	0.737	15.50	1.10	0.68	0.096
19.20	1.10	1.10	0.219	1.90	1.82	4.35	0.729	16.00	1.08	0.66	0.092
19.60	1.05	1.11	0.227	2.00	1.75	4.18	0.721	16.50	1.06	0.63	0.086
20.00	0.96	1.10	0.239	2.10	1.67	4.03	0.714	17.00	1.07	0.61	0.081
20.40	0.93	1.09	0.240	2.20	1.60	3.88	0.707	17.50	1.06	0.61	0.080
20.80	0.89	1.05	0.240	2.30	1.53	3.75	0.700	18.00	1.07	0.59	0.077
21.20	0.86	1.02	0.237	2.40	1.47	3.61	0.693	18.50	1.07	0.59	0.077
21.60	0.83	0.99	0.235	2.50	1.41	3.48	0.685	19.00	1.08	0.59	0.077
22.00	0.80	0.96	0.230	2.60	1.37	3.36	0.676	19.50	1.08	0.61	0.080
22.40	0.78	0.93	0.226	2.70	1.32	3.25	0.668	20.00	1.07	0.65	0.090
22.80	0.77	0.90	0.220	2.80	1.29	3.13	0.658	20.50	1.03	0.67	0.098
23.20	0.75	0.88	0.217	2.90	1.26	3.03	0.648	21.00	0.99	0.67	0.103
23.60	0.75	0.86	0.211	3.00	1.23	2.94	0.639	21.50	0.95	0.66	0.103
24.00	0.73	0.84	0.209	3.10	1.20	2.85	0.630	22.00	0.91	0.64	0.103
24.40	0.72	0.82	0.207	3.20	1.17	2.77	0.622	22.50	0.88	0.62	0.101
24.80	0.70	0.80	0.205	3.30	1.14	2.68	0.613	23.00	0.86	0.59	0.097
25.20	0.69	0.77	0.202	3.40	1.12	2.60	0.602	23.50	0.85	0.56	0.091
25.60	0.67	0.75	0.199	3.50	1.10	2.52	0.591	24.00	0.84	0.54	0.086
26.00	0.66	0.72	0.195	3.60	1.08	2.45	0.581	25.00	0.81	0.51	0.084
26.40	0.65	0.69	0.189	3.70	1.07	2.38	0.570	26.40	0.80	0.43	0.066
26.80	0.63	0.66	0.183	3.80	1.06	2.31	0.558	27.80	0.81	0.38	0.052
27.20	0.65	0.62	0.165	3.90	1.05	2.25	0.547	29.20	0.82	0.35	0.046
28.00	0.64	0.59	0.156	4.00	1.03	2.19	0.537				
28.40	0.64	0.57	0.148	4.20	1.04	2.09	0.510				
28.80	0.65	0.55	0.140	4.40	1.03	2.01	0.493	<i>Platinum</i> ²⁰			
29.20	0.65	0.53	0.134	4.60	1.03	1.94	0.476	0.10	13.21	44.72	0.976
29.60	0.65	0.51	0.128	4.80	1.01	1.90	0.470	0.15	8.18	31.16	0.969
30.00	0.65	0.49	0.121	5.00	0.96	1.86	0.472	0.20	5.90	23.95	0.962
31.00	0.65	0.45	0.111	5.20	0.90	1.79	0.474	0.25	4.70	19.40	0.954
32.00	0.66	0.41	0.095	5.40	0.85	1.70	0.463	0.30	3.92	16.16	0.945
33.00	0.68	0.37	0.079	5.60	0.81	1.62	0.449	0.35	3.28	13.66	0.936
34.00	0.70	0.34	0.068	5.80	0.78	1.54	0.437	0.40	2.81	11.38	0.922
35.00	0.72	0.31	0.057	6.00	0.76	1.45	0.418	0.45	3.03	9.31	0.882
36.00	0.74	0.29	0.048	6.20	0.74	1.37	0.397	0.50	3.91	7.71	0.813
37.00	0.77	0.27	0.040	6.40	0.73	1.29	0.375	0.55	4.58	7.14	0.777
38.00	0.79	0.26	0.035	6.60	0.72	1.21	0.350	0.60	5.13	6.75	0.753
39.00	0.81	0.26	0.031	6.80	0.73	1.13	0.316	0.65	5.52	6.66	0.746
40.00	0.84	0.26	0.026	7.00	0.73	1.05	0.287	0.70	5.71	6.83	0.751
				7.20	0.75	0.98	0.255	0.75	5.57	7.02	0.759
				7.40	0.77	0.91	0.223	0.80	5.31	7.04	0.762
				7.60	0.79	0.85	0.195	0.85	5.05	6.98	0.763
				7.80	0.83	0.78	0.163	0.90	4.77	6.91	0.765
				8.00	0.88	0.73	0.133	0.95	4.50	6.77	0.763
				8.20	0.94	0.70	0.117	1.00	4.25	6.62	0.762
				8.40	0.96	0.70	0.114	1.10	3.86	6.24	0.753
				8.60	1.00	0.65	0.097	1.20	3.55	5.92	0.746
				8.80	1.04	0.65	0.094	1.30	3.29	5.61	0.736
				9.00	1.07	0.64	0.090	1.40	3.10	5.32	0.725
								1.50	2.92	5.07	0.716
								1.60	2.76	4.84	0.706

Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$
1.70	2.63	4.64	0.697	12.80	1.29	0.94	0.158	4.00	0.34	0.08	0.245
1.80	2.51	4.43	0.686	13.20	1.31	0.93	0.155	4.065	0.38	0.07	0.204
1.90	2.38	4.26	0.678	13.60	1.31	0.93	0.155	4.133	0.41	0.07	0.177
2.00	2.30	4.07	0.664	14.00	1.31	0.93	0.155	4.203	0.45	0.06	0.145
2.10	2.23	3.92	0.654	14.40	1.30	0.93	0.156	4.275	0.48	0.06	0.125
2.20	2.17	3.77	0.642	14.80	1.27	0.93	0.157	4.350	0.52	0.05	0.101
2.30	2.10	3.67	0.636	15.20	1.27	0.93	0.155	4.428	0.55	0.05	0.085
2.40	2.03	3.54	0.626	15.60	1.25	0.92	0.151	4.509	0.58	0.05	0.072
2.50	1.96	3.42	0.616	16.00	1.24	0.89	0.146	4.592	0.61	0.05	0.060
2.60	1.91	3.30	0.605	16.50	1.24	0.87	0.142	4.679	0.64	0.04	0.049
2.70	1.87	3.20	0.595	17.00	1.25	0.86	0.138	4.769	0.66	0.04	0.043
2.80	1.83	3.10	0.585	17.50	1.27	0.85	0.135	4.862	0.68	0.04	0.037
2.90	1.79	3.01	0.575	18.00	1.31	0.88	0.142	4.959	0.70	0.04	0.032
3.00	1.75	2.92	0.565	18.50	1.30	0.94	0.157	5.061	0.72	0.04	0.027
3.20	1.68	2.76	0.546	19.00	1.28	0.99	0.171	5.166	0.74	0.04	0.023
3.40	1.63	2.62	0.527	19.50	1.23	1.03	0.184	5.276	0.76	0.04	0.019
3.60	1.58	2.48	0.507	20.00	1.18	1.06	0.197	5.391	0.78	0.04	0.016
3.80	1.53	2.37	0.491	20.50	1.11	1.09	0.212	5.510	0.79	0.05	0.015
4.00	1.49	2.25	0.472	21.00	1.03	1.10	0.226	5.637	0.81	0.05	0.012
4.20	1.45	2.14	0.452	21.50	0.94	1.08	0.238	5.767	0.83	0.05	0.009
4.40	1.43	2.04	0.432	22.00	0.87	1.04	0.240	6.048	0.85	0.05	0.007
4.60	1.39	1.95	0.415	22.50	0.81	0.98	0.235	6.199	0.87	0.05	0.006
4.80	1.38	1.85	0.392	23.00	0.77	0.92	0.226	6.358	0.88	0.05	0.005
5.00	1.36	1.76	0.372	23.50	0.75	0.87	0.213	6.526	0.90	0.06	0.004
5.20	1.36	1.67	0.350	24.00	0.74	0.82	0.201	6.702	0.91	0.06	0.003
5.40	1.36	1.61	0.332	24.50	0.73	0.77	0.187	6.888	0.92	0.06	0.003
5.60	1.36	1.54	0.315	25.00	0.73	0.73	0.174	7.085	0.92	0.06	0.003
5.80	1.36	1.47	0.295	25.50	0.73	0.70	0.162	7.293	0.93	0.06	0.002
6.00	1.38	1.40	0.276	26.00	0.74	0.67	0.150	7.514	0.93	0.06	0.002
6.20	1.39	1.35	0.261	26.50	0.74	0.65	0.142	7.749	0.94	0.06	0.002
6.40	1.42	1.29	0.246	27.00	0.74	0.63	0.136	7.999	0.94	0.06	0.002
6.60	1.45	1.26	0.236	27.50	0.74	0.62	0.130	8.260	0.94	0.06	0.002
6.80	1.48	1.24	0.231	28.00	0.75	0.60	0.125	8.551	0.94	0.06	0.002
7.00	1.50	1.24	0.230	28.50	0.75	0.59	0.121	8.856	0.94	0.05	0.002
7.20	1.50	1.25	0.231	29.00	0.75	0.58	0.118	9.184	0.94	0.05	0.002
7.40	1.49	1.23	0.228	29.50	0.74	0.58	0.120	9.537	0.94	0.04	0.001
7.60	1.48	1.22	0.225	30.00	0.73	0.58	0.124	9.919	0.94	0.04	0.001
7.80	1.48	1.20	0.221					10.33	0.94	0.03	0.001
8.00	1.47	1.18	0.216	<i>Potassium</i> ²¹				11.0		0.03	
8.20	1.47	1.17	0.212	0.55	0.139	7.10	0.989	12.0		0.028	
8.40	1.47	1.15	0.209	0.58	0.119	6.72	0.990				
8.60	1.47	1.14	0.205	0.63	0.106	6.32	0.990	<i>Rhenium, single crystal, $\bar{E} \parallel c^p$</i>			
8.80	1.47	1.13	0.202	0.67	0.091	5.79	0.990	0.10	6.06	51.03	0.991
9.00	1.48	1.12	0.200	0.73	0.079	5.30	0.989	0.15	4.66	33.96	0.984
9.20	1.49	1.11	0.198	0.81	0.066	4.75	0.989	0.20	4.16	25.36	0.975
9.40	1.49	1.12	0.200	0.92	0.056	4.19	0.988	0.25	4.03	20.10	0.962
9.60	1.49	1.13	0.203	1.05	0.044	3.58	0.987	0.30	4.37	16.69	0.943
9.80	1.48	1.15	0.207	1.23	0.040	3.04	0.985	0.35	4.50	14.53	0.925
10.00	1.46	1.15	0.209	1.44	0.040	2.56	0.979	0.40	4.53	12.96	0.909
10.20	1.43	1.16	0.211	1.65	0.044	2.19	0.970	0.45	4.53	11.78	0.893
10.40	1.40	1.15	0.210	1.87	0.050	1.84	0.955	0.50	4.53	10.88	0.878
10.60	1.37	1.14	0.207	2.07	0.053	1.62	0.943	0.55	4.50	10.26	0.867
10.80	1.35	1.12	0.203	2.27	0.049	1.43	0.938	0.60	4.29	9.75	0.861
11.00	1.33	1.10	0.199	2.45	0.046	1.28	0.933	0.65	4.07	9.35	0.856
11.20	1.31	1.08	0.194	2.64	0.043	1.14	0.928	0.70	3.80	8.94	0.853
11.40	1.30	1.06	0.188	2.82	0.043	1.02	0.919	0.75	3.48	8.55	0.850
11.60	1.29	1.04	0.183	2.95	0.041	0.898	0.913	0.80	3.21	8.10	0.846
11.80	1.29	1.01	0.177	3.06	0.041	0.799	0.905	0.85	2.96	7.68	0.841
12.00	1.29	1.00	0.173	3.40	0.052	0.549	0.852	0.90	2.73	7.24	0.835
12.40	1.29	0.97	0.165	3.71	0.089	0.288	0.719	0.95	2.56	6.79	0.826
				3.97	0.287	0.091	0.310	1.00	2.45	6.36	0.813

Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$
1.10	2.38	5.61	0.778	11.40	1.28	1.28	0.252	50.00	0.80	0.30	0.038
1.20	2.35	5.02	0.742	11.60	1.26	1.28	0.252	52.00	0.78	0.30	0.044
1.30	2.39	4.54	0.702	11.80	1.24	1.26	0.249	54.00	0.72	0.30	0.055
1.40	2.44	4.13	0.662	12.00	1.23	1.24	0.244	56.00	0.66	0.24	0.061
1.50	2.50	3.79	0.624	12.40	1.22	1.21	0.237	58.00	0.65	0.16	0.055
1.60	2.59	3.49	0.587	12.80	1.21	1.18	0.230	<i>Rhenium, single crystal, $\bar{E} \perp \bar{c}^0$</i>			
1.70	2.70	3.27	0.557	13.20	1.22	1.16	0.222	0.10	4.25	42.83	0.991
1.80	2.82	3.10	0.535	13.60	1.22	1.13	0.215	0.15	3.28	28.08	0.984
1.90	2.90	3.00	0.520	14.00	1.24	1.12	0.209	0.20	3.28	20.66	0.971
2.00	2.97	2.91	0.510	14.40	1.27	1.11	0.204	0.25	3.47	16.27	0.951
2.10	3.03	2.86	0.504	14.80	1.29	1.15	0.213	0.30	3.73	13.44	0.926
2.20	3.06	2.84	0.501	15.20	1.29	1.19	0.225	0.35	3.93	11.54	0.900
2.30	3.07	2.82	0.499	15.60	1.26	1.22	0.236	0.40	3.99	10.15	0.875
2.40	3.06	2.81	0.498	16.00	1.23	1.25	0.248	0.45	4.17	9.03	0.846
2.50	3.02	2.80	0.497	16.40	1.19	1.27	0.259	0.50	4.34	8.26	0.821
2.60	2.96	2.77	0.493	16.80	1.14	1.29	0.269	0.55	4.45	7.73	0.801
2.70	2.89	2.68	0.482	17.00	1.12	1.30	0.275	0.60	4.53	7.40	0.788
2.80	2.89	2.57	0.468	17.40	1.07	1.30	0.286	0.65	4.44	7.26	0.784
2.90	2.99	2.47	0.457	18.00	0.99	1.30	0.300	0.70	4.13	7.09	0.784
3.00	3.11	2.57	0.470	18.40	0.93	1.29	0.311	0.75	3.77	6.75	0.779
3.20	2.90	2.68	0.482	18.80	0.87	1.28	0.321	0.80	3.55	6.32	0.766
3.40	2.83	2.50	0.459	19.20	0.81	1.25	0.330	0.85	3.39	5.95	0.752
3.60	2.93	2.48	0.457	19.60	0.77	1.21	0.332	0.90	3.26	5.61	0.737
3.80	2.86	2.56	0.467	20.00	0.73	1.18	0.333	0.95	3.17	5.27	0.719
4.00	2.81	2.51	0.460	20.40	0.70	1.14	0.332	1.00	3.09	4.96	0.701
4.20	2.86	2.55	0.466	20.80	0.67	1.11	0.332	1.10	3.05	4.39	0.658
4.40	2.81	2.74	0.489	21.20	0.64	1.08	0.334	1.20	3.08	3.89	0.613
4.60	2.56	2.83	0.504	21.60	0.61	1.04	0.335	1.30	3.20	3.56	0.578
4.80	2.41	2.71	0.493	22.00	0.58	1.01	0.340	1.40	3.23	3.38	0.559
5.00	2.39	2.68	0.488	22.40	0.55	0.97	0.341	1.50	3.23	3.12	0.532
5.20	2.34	2.75	0.500	22.80	0.53	0.93	0.338	1.60	3.29	2.88	0.507
5.40	2.20	2.81	0.515	23.20	0.51	0.89	0.334	1.70	3.38	2.72	0.491
5.60	2.02	2.84	0.530	23.60	0.50	0.85	0.329	1.80	3.47	2.59	0.480
5.80	1.83	2.80	0.538	24.00	0.48	0.80	0.319	1.90	3.54	2.50	0.473
6.00	1.65	2.71	0.541	24.40	0.48	0.76	0.207	2.00	3.63	2.43	0.469
6.20	1.54	2.59	0.532	24.80	0.47	0.72	0.296	2.10	3.74	2.40	0.470
6.40	1.45	2.50	0.526	25.20	0.47	0.68	0.282	2.20	3.83	2.38	0.472
6.80	1.32	2.31	0.508	25.60	0.47	0.65	0.270	2.30	3.93	2.44	0.481
7.00	1.26	2.23	0.500	26.00	0.47	0.61	0.255	2.40	4.00	2.55	0.492
7.20	1.20	2.15	0.493	26.40	0.48	0.57	0.240	2.50	4.01	2.70	0.505
7.40	1.16	2.06	0.480	26.80	0.48	0.54	0.225	2.60	3.90	2.84	0.514
7.60	1.12	1.99	0.470	27.20	0.49	0.51	0.208	2.70	3.74	2.92	0.517
7.80	1.08	1.89	0.454	27.60	0.50	0.48	0.193	2.80	3.57	2.88	0.511
8.00	1.05	1.80	0.435	28.00	0.51	0.45	0.176	2.90	3.49	2.75	0.497
8.20	1.05	1.71	0.411	29.00	0.54	0.39	0.145	3.00	3.53	2.71	0.493
8.40	1.05	1.62	0.386	30.00	0.57	0.33	0.114	3.20	3.55	2.84	0.506
8.60	1.06	1.55	0.360	31.00	0.62	0.29	0.086	3.40	3.34	2.88	0.508
8.80	1.09	1.48	0.336	32.00	0.66	0.26	0.065	3.60	3.25	2.83	0.501
9.00	1.11	1.43	0.317	33.00	0.68	0.24	0.054	3.80	3.24	2.84	0.502
9.20	1.13	1.39	0.301	34.00	0.72	0.21	0.041	4.00	3.19	2.94	0.513
9.40	1.16	1.34	0.281	35.00	0.76	0.20	0.031	4.20	3.05	3.06	0.526
9.60	1.18	1.32	0.274	36.00	0.79	0.20	0.025	4.40	2.88	3.15	0.539
9.80	1.20	1.29	0.264	37.00	0.82	0.19	0.021	4.60	2.67	3.18	0.548
10.00	1.23	1.26	0.252	38.00	0.85	0.20	0.018	4.80	2.44	3.17	0.554
10.20	1.25	1.25	0.246	39.00	0.89	0.21	0.016	5.00	2.25	3.12	0.556
10.40	1.28	1.25	0.242	40.00	0.88	0.26	0.022	5.20	2.10	3.04	0.555
10.60	1.29	1.25	0.242	42.00	0.88	0.26	0.022	5.40	1.96	2.96	0.553
10.80	1.30	1.26	0.244	44.00	0.89	0.29	0.026	5.60	1.84	2.88	0.551
11.00	1.30	1.27	0.247	46.00	0.85	0.32	0.035	5.80	1.73	2.81	0.549
11.20	1.29	1.28	0.249	48.00	0.82	0.30	0.036	6.00	1.61	2.74	0.549

Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$
6.20	1.51	2.64	0.545	24.80	0.48	0.75	0.303	3.10	1.41	4.20	0.760
6.40	1.42	2.56	0.541	25.20	0.47	0.72	0.295	3.20	1.30	4.09	0.764
6.80	1.28	2.37	0.526	25.60	0.47	0.68	0.286	3.30	1.20	3.97	0.767
7.00	1.22	2.28	0.517	26.00	0.46	0.64	0.276	3.40	1.11	3.84	0.769
7.20	1.16	2.19	0.508	26.40	0.46	0.61	0.263	3.50	1.04	3.71	0.768
7.40	1.12	2.08	0.493	26.80	0.46	0.57	0.249	3.60	0.99	3.58	0.764
7.60	1.12	1.98	0.468	27.20	0.47	0.53	0.231	3.70	0.95	3.45	0.759
7.80	1.08	1.93	0.463	27.60	0.48	0.50	0.216	3.80	0.91	3.34	0.753
8.00	1.05	1.83	0.443	28.00	0.49	0.47	0.198	3.90	0.88	3.23	0.747
8.20	1.05	1.74	0.418	29.00	0.51	0.41	0.164	4.00	0.86	3.12	0.739
8.40	1.05	1.66	0.397	30.00	0.55	0.34	0.129	4.20	0.83	2.94	0.722
8.60	1.06	1.58	0.372	31.00	0.59	0.29	0.097	4.40	0.80	2.76	0.706
8.80	1.07	1.52	0.351	32.00	0.64	0.26	0.072	4.60	0.78	2.60	0.684
9.00	1.09	1.46	0.327	33.00	0.67	0.24	0.060	4.80	0.79	2.46	0.659
9.20	1.11	1.41	0.309	34.00	0.70	0.22	0.047	5.00	0.79	2.34	0.635
9.40	1.14	1.36	0.290	35.00	0.74	0.20	0.036	5.20	0.79	2.23	0.613
9.60	1.17	1.31	0.273	36.00	0.77	0.19	0.029	5.40	0.80	2.14	0.591
9.80	1.20	1.27	0.258	37.00	0.80	0.19	0.023	5.60	0.80	2.06	0.573
10.00	1.24	1.24	0.244	38.00	0.84	0.19	0.018	5.80	0.79	2.00	0.561
10.20	1.29	1.22	0.234	39.00	0.88	0.21	0.016	6.00	0.76	1.93	0.556
10.40	1.33	1.23	0.233	40.00	0.87	0.25	0.023	6.20	0.73	1.85	0.544
10.60	1.36	1.25	0.238	42.00	0.87	0.25	0.023	6.40	0.70	1.77	0.534
10.80	1.38	1.28	0.245	44.00	0.88	0.28	0.026	6.60	0.68	1.69	0.518
11.00	1.37	1.31	0.253	46.00	0.84	0.31	0.035	6.80	0.67	1.60	0.498
11.20	1.36	1.33	0.259	48.00	0.82	0.30	0.036	7.00	0.66	1.52	0.476
11.40	1.33	1.34	0.264	50.00	0.80	0.30	0.039	7.20	0.66	1.43	0.452
11.60	1.31	1.34	0.266	52.00	0.77	0.30	0.044	7.40	0.66	1.35	0.423
11.80	1.28	1.33	0.266	54.00	0.71	0.29	0.055	7.60	0.67	1.27	0.394
12.00	1.26	1.32	0.264	56.00	0.66	0.23	0.061	7.80	0.68	1.20	0.363
12.40	1.23	1.29	0.257	58.00	0.64	0.16	0.055	8.00	0.69	1.12	0.329
12.80	1.22	1.26	0.251					8.20	0.71	1.04	0.288
13.20	1.20	1.23	0.245	<i>Rhodium^{II}</i>				8.40	0.74	0.97	0.252
13.60	1.19	1.20	0.236	0.10	18.48	69.43	0.986	8.60	0.78	0.89	0.212
14.00	1.20	1.16	0.225	0.20	8.66	37.46	0.977	8.80	0.83	0.83	0.179
14.40	1.22	1.13	0.214	0.30	5.85	25.94	0.967	9.00	0.88	0.77	0.148
14.80	1.27	1.12	0.207	0.40	4.74	19.80	0.955	9.20	0.95	0.73	0.125
15.20	1.31	1.17	0.218	0.50	4.20	16.07	0.941	9.40	1.01	0.71	0.110
15.60	1.31	1.23	0.234	0.60	3.87	13.51	0.925	9.60	1.07	0.69	0.102
16.00	1.28	1.28	0.251	0.70	3.67	11.72	0.908	9.80	1.12	0.69	0.098
16.40	1.24	1.33	0.270	0.80	3.63	10.34	0.887	10.00	1.17	0.69	0.098
16.80	1.17	1.37	0.288	0.90	3.62	9.36	0.867	10.60	1.26	0.73	0.106
17.00	1.14	1.38	0.297	1.00	3.71	8.67	0.848	11.00	1.29	0.76	0.113
17.40	1.06	1.39	0.314	1.10	3.67	8.26	0.837	11.60	1.32	0.80	0.124
18.00	0.95	1.38	0.334	1.20	3.51	7.94	0.832	12.00	1.32	0.82	0.127
18.40	0.88	1.36	0.346	1.30	3.26	7.63	0.829	12.60	1.32	0.82	0.129
18.80	0.82	1.33	0.355	1.40	3.01	7.31	0.827	13.00	1.32	0.83	0.131
19.20	0.76	1.29	0.360	1.50	2.78	6.97	0.823	13.60	1.32	0.85	0.134
19.60	0.72	1.25	0.363	1.60	2.60	6.64	0.818	14.00	1.32	0.86	0.138
20.00	0.67	1.21	0.369	1.70	2.42	6.33	0.813	14.60	1.30	0.89	0.144
20.40	0.64	1.15	0.364	1.80	2.30	6.02	0.805	15.00	1.28	0.90	0.147
20.80	0.61	1.10	0.357	1.90	2.20	5.76	0.798	15.60	1.25	0.90	0.147
21.20	0.60	1.06	0.349	2.00	2.12	5.51	0.789	16.00	1.24	0.89	0.147
21.60	0.58	1.02	0.342	2.10	2.05	5.30	0.780	16.50	1.23	0.88	0.145
22.00	0.57	0.98	0.336	2.20	2.00	5.11	0.772	17.00	1.22	0.88	0.144
22.40	0.56	0.95	0.328	2.30	1.94	4.94	0.765	17.50	1.22	0.87	0.143
22.80	0.55	0.92	0.325	2.40	1.90	4.78	0.756	18.00	1.23	0.88	0.145
23.20	0.53	0.89	0.322	2.50	1.88	4.65	0.748	18.50	1.25	0.92	0.155
23.60	0.52	0.85	0.317	2.60	1.85	4.55	0.743	19.00	1.24	0.98	0.172
24.00	0.50	0.82	0.314	2.70	1.80	4.49	0.742	19.50	1.18	1.05	0.193
24.40	0.49	0.79	0.309	2.90	1.63	4.36	0.748	20.00	1.10	1.09	0.213
				3.00	1.53	4.29	0.753				

Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$
20.50	1.00	1.09	0.230	3.30	2.00	3.91	0.671	17.50	1.32	0.93	0.155
21.00	0.91	1.05	0.234	3.40	1.87	3.83	0.673	18.00	1.26	0.99	0.173
21.50	0.86	1.00	0.228	3.50	1.76	3.74	0.674	18.50	1.18	1.02	0.185
22.00	0.83	0.95	0.219	3.60	1.66	3.65	0.675	19.00	1.11	1.02	0.192
22.50	0.81	0.92	0.214	3.70	1.57	3.55	0.673	19.50	1.05	1.02	0.199
23.00	0.79	0.90	0.213	3.80	1.49	3.45	0.672	20.00	0.99	1.02	0.208
23.50	0.75	0.87	0.214	3.90	1.42	3.35	0.668	20.50	0.92	0.99	0.212
24.00	0.73	0.84	0.210	4.00	1.37	3.24	0.661	21.00	0.86	0.94	0.209
24.50	0.70	0.81	0.208	4.20	1.29	3.08	0.649	21.50	0.83	0.90	0.203
25.00	0.69	0.77	0.202	4.40	1.22	2.93	0.639	22.00	0.81	0.86	0.193
25.50	0.67	0.74	0.195	4.60	1.16	2.79	0.628	23.00	0.77	0.79	0.182
26.00	0.66	0.70	0.188	4.80	1.11	2.67	0.617	24.00	0.74	0.74	0.171
26.50	0.65	0.66	0.176	5.00	1.06	2.56	0.607	25.00	0.71	0.69	0.163
27.00	0.65	0.64	0.168	5.20	1.01	2.46	0.600	26.00	0.68	0.63	0.154
27.50	0.65	0.61	0.159	5.40	0.95	2.35	0.593	27.00	0.67	0.57	0.140
28.00	0.65	0.59	0.152	5.60	0.92	2.23	0.576	28.00	0.66	0.51	0.124
29.00	0.65	0.54	0.137	5.80	0.90	2.14	0.559	29.00	0.67	0.46	0.107
30.00	0.66	0.51	0.127	6.00	0.88	2.05	0.545	30.00	0.67	0.43	0.097
31.00	0.64	0.49	0.127	6.20	0.87	1.98	0.531	31.00	0.67	0.37	0.084
32.00	0.61	0.44	0.126	6.40	0.84	1.91	0.521	32.00	0.69	0.33	0.070
33.00	0.60	0.37	0.110	6.60	0.82	1.84	0.510	33.00	0.71	0.30	0.058
34.00	0.65	0.30	0.074	6.80	0.79	1.77	0.500	34.00	0.73	0.27	0.048
35.00	0.69	0.28	0.058	7.00	0.76	1.69	0.489	35.00	0.75	0.25	0.039
36.00	0.73	0.27	0.049	7.20	0.75	1.61	0.472	36.00	0.77	0.24	0.035
37.00	0.74	0.28	0.047	7.40	0.73	1.54	0.455	37.00	0.79	0.23	0.039
38.00	0.74	0.27	0.045	7.60	0.73	1.46	0.433	38.00	0.80	0.22	0.027
39.00	0.75	0.25	0.041	7.80	0.73	1.39	0.411	39.00	0.82	0.22	0.024
<i>Ruthenium, single crystal, $\vec{E} \parallel \vec{c}^o$</i>				8.00	0.72	1.33	0.391	40.00	0.83	0.22	0.022
0.10	11.50	51.38	0.984	8.20	0.72	1.26	0.366	<i>Ruthenium, single crystal, $\vec{E} \perp \vec{c}^o$</i>			
0.20	5.93	27.14	0.970	8.40	0.73	1.20	0.342	0.10	11.85	50.81	0.983
0.30	4.33	18.50	0.953	8.60	0.74	1.14	0.318	0.20	6.68	27.18	0.966
0.40	3.60	13.97	0.933	8.80	0.74	1.08	0.295	0.30	4.94	18.92	0.950
0.50	3.18	11.04	0.909	9.00	0.75	1.02	0.267	0.40	3.90	14.51	0.933
0.60	3.28	8.89	0.865	9.20	0.77	0.97	0.243	0.50	3.27	11.63	0.915
0.70	3.62	7.73	0.822	9.40	0.79	0.91	0.217	0.60	2.98	9.54	0.888
0.80	3.42	7.02	0.801	9.60	0.82	0.86	0.190	0.70	2.82	7.99	0.856
0.90	3.25	6.12	0.766	9.80	0.85	0.81	0.167	0.80	2.73	6.71	0.815
1.00	3.39	5.33	0.715	10.00	0.88	0.76	0.144	0.90	2.82	5.54	0.751
1.10	3.66	4.83	0.675	10.20	0.92	0.72	0.125	1.00	3.17	4.59	0.670
1.20	3.84	4.57	0.654	10.40	0.96	0.69	0.110	1.10	3.69	3.91	0.604
1.30	3.94	4.38	0.638	10.60	1.01	0.67	0.100	1.20	4.28	3.66	0.585
1.40	4.02	4.19	0.624	10.80	1.05	0.66	0.094	1.30	4.66	3.72	0.593
1.50	4.16	4.07	0.614	11.00	1.09	0.65	0.090	1.40	4.86	3.79	0.601
1.60	4.33	4.08	0.615	11.20	1.12	0.65	0.088	1.50	4.99	3.89	0.609
1.70	4.42	4.21	0.624	11.40	1.15	0.65	0.087	1.60	5.08	4.03	0.618
1.80	4.40	4.38	0.636	11.60	1.18	0.65	0.088	1.70	5.12	4.22	0.629
1.90	4.29	4.61	0.651	11.80	1.21	0.66	0.090	1.80	5.10	4.45	0.642
2.00	4.04	4.81	0.667	12.00	1.23	0.67	0.092	1.90	4.96	4.78	0.660
2.10	3.69	4.90	0.679	12.40	1.26	0.69	0.098	2.00	4.61	5.06	0.677
2.20	3.35	4.82	0.683	12.80	1.27	0.72	0.104	2.10	4.21	5.09	0.682
2.30	3.09	4.70	0.681	13.20	1.28	0.74	0.108	2.20	3.94	5.00	0.681
2.40	2.89	4.55	0.677	13.60	1.28	0.75	0.111	2.30	3.69	4.97	0.684
2.50	2.74	4.40	0.671	14.00	1.28	0.76	0.114	2.40	3.44	4.88	0.684
2.60	2.64	4.25	0.663	14.40	1.27	0.76	0.114	2.50	3.27	4.77	0.681
2.70	2.58	4.14	0.656	14.80	1.27	0.76	0.114	2.60	3.14	4.66	0.677
2.80	2.54	4.05	0.650	15.00	1.27	0.76	0.114	2.70	3.06	4.59	0.674
2.90	2.48	4.03	0.650	15.60	1.28	0.77	0.115	2.80	2.99	4.59	0.676
3.00	2.38	4.03	0.656	16.00	1.30	0.78	0.118	2.90	2.87	4.64	0.686
3.10	2.26	4.00	0.661	16.50	1.32	0.80	0.123	3.00	2.64	4.69	0.701
3.20	2.13	3.96	0.666	17.00	1.34	0.85	0.136	3.10	2.40	4.64	0.710

Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)	Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)	Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)		
3.20	2.18	4.55	0.717	17.00	1.28	0.94	0.158	0.4959	3.442	1.41E-04	0.302		
3.30	2.00	4.43	0.721	17.50	1.25	1.00	0.175	0.6199	3.462	1.12E-04	0.304		
3.40	1.84	4.30	0.723	18.00	1.19	1.04	0.190	0.7439	3.486	9.42E-05	0.307		
3.50	1.71	4.16	0.723	18.50	1.12	1.05	0.200	0.8679	3.516	8.07E-05	0.310		
3.60	1.60	4.03	0.722	19.00	1.07	1.05	0.205	0.9919	3.551	7.11E-05	0.314		
3.70	1.50	3.90	0.721	19.50	1.02	1.04	0.212	1.116	3.592	6.37E-05	0.319		
3.80	1.41	3.77	0.718	20.00	0.97	1.04	0.219	1.240	3.640	5.81E-05	0.324		
3.90	1.35	3.64	0.713	20.50	0.91	1.03	0.228	1.50		1.33E-04			
4.00	1.29	3.53	0.707	21.00	0.85	1.01	0.234	1.60		1.59E-04			
4.20	1.21	3.31	0.694	21.50	0.80	0.97	0.234	1.70		6.27E-04			
4.40	1.16	3.13	0.679	22.00	0.77	0.94	0.233	1.80	4.46	2.20E-02	0.402		
4.60	1.13	2.97	0.662	23.00	0.71	0.87	0.229	2.0	4.79	0.76	0.438		
4.80	1.09	2.86	0.652	24.00	0.67	0.79	0.218	2.2	4.49	1.19	0.431		
5.00	1.03	2.75	0.648	25.00	0.64	0.73	0.205	2.4	4.28	1.21	0.417		
5.20	0.97	2.64	0.643	26.00	0.61	0.66	0.194	2.6	4.40	1.32	0.430		
5.40	0.91	2.52	0.635	27.00	0.60	0.59	0.177	2.8	4.59	1.70	0.462		
5.60	0.88	2.40	0.622	28.00	0.60	0.53	0.155	3.0	4.44	2.29	0.490		
5.80	0.86	2.29	0.605	29.00	0.61	0.48	0.134	3.2	3.92	2.59	0.493		
6.00	0.84	2.20	0.591	30.00	0.62	0.45	0.123	3.4	3.69	2.76	0.502		
6.20	0.82	2.11	0.576	31.00	0.61	0.40	0.114	3.6	3.39	3.01	0.521		
6.40	0.81	2.04	0.564	32.00	0.63	0.34	0.093	3.8	(3.00)				
6.60	0.78	1.97	0.556	33.00	0.65	0.31	0.077	4.0	(2.65)				
6.80	0.76	1.89	0.545	34.00	0.67	0.28	0.065	4.2	(2.30)				
7.00	0.73	1.82	0.538	35.00	0.70	0.26	0.054	4.5	1.92	2.78	0.528		
7.20	0.70	1.75	0.527	36.00	0.72	0.25	0.047	5.0	1.50	2.31	0.482		
7.40	0.68	1.67	0.513	37.00	0.73	0.23	0.041	6.0	1.57	1.49	0.288		
7.60	0.67	1.59	0.496	38.00	0.75	0.22	0.035	7.0	1.84	1.45	0.276		
7.80	0.66	1.51	0.476	39.00	0.77	0.22	0.031	8.0	1.35	1.68	0.353		
8.00	0.66	1.44	0.454	40.00	0.79	0.22	0.028	9.0	1.35	1.64	0.342		
8.20	0.65	1.36	0.430					10.0	0.92	1.07	0.238		
8.40	0.66	1.29	0.403					12.0	1.00	1.10	0.232		
8.60	0.66	1.22	0.378	<i>Selenium, single crystal, $\vec{E} \parallel \hat{c}^{22}$</i>	0.01364	2.914	0.248	0.242	14.0	0.81	0.91	0.211	
8.80	0.68	1.15	0.346		0.01488	3.175	9.95E-02	0.272	16.0	0.65	0.61	0.160	
9.00	0.69	1.09	0.317		0.01612	3.263	2.13E-03	0.282	18.0	0.65	0.48	0.120	
9.20	0.70	1.02	0.286		0.01736	3.306	3.81E-02	0.287	20.0	0.69	0.36	0.076	
9.40	0.73	0.95	0.251		0.01860	3.330	7.04E-03	0.290	22.0	0.81	0.25	0.030	
9.60	0.77	0.89	0.216		0.01984	3.346	4.23E-02	0.291	24.0	0.91	0.18	0.011	
9.80	0.82	0.84	0.185		0.02108	3.358	3.40E-03	0.293	26.0	0.86	0.15	0.012	
10.00	0.86	0.81	0.163		0.02232	3.366	5.31E-02	0.294	28.0	0.85	0.13	0.011	
10.20	0.90	0.77	0.143		0.02356	3.372	1.96E-03	0.294	30.0	0.87	0.11	0.008	
10.40	0.94	0.74	0.127		0.02480	3.377	2.39E-02	0.295					
10.60	0.99	0.72	0.115		0.02604	3.380		0.295	<i>Selenium, single crystal, $\vec{E} \perp \hat{c}^{22}$</i>	0.01364	2.854	0.0239	0.231
10.80	1.04	0.71	0.108		0.02728		1.16E-02			0.01488	2.932	0.0325	0.241
11.00	1.08	0.70	0.104		0.02976		7.96E-03			0.01612	3.140	0.1750	0.269
11.20	1.11	0.70	0.102		0.03224		8.57E-03			0.01736	2.959	1.3300	0.321
11.40	1.14	0.70	0.101		0.03472		2.70E-02			0.01860	2.111	0.2550	0.133
11.60	1.17	0.71	0.102		0.03720	3.397	1.72E-02	0.297		0.01984	2.356	0.0746	0.164
11.80	1.20	0.72	0.104		0.04463		1.13E-02			0.02108	2.462	0.0276	0.178
12.00	1.22	0.73	0.107		0.04959	3.403	2.79E-03	0.298		0.02232	2.502	0.0442	0.184
12.40	1.25	0.76	0.113		0.05703		1.56E-03			0.02356	2.543	0.0097	0.190
12.80	1.26	0.78	0.118		0.06199	3.405	1.35E-03	0.298		0.02480	2.550	0.0239	0.191
13.20	1.27	0.81	0.124		0.06819		5.79E-04			0.02604	2.582		0.195
13.60	1.27	0.83	0.129		0.07439	3.407	4.44E-04	0.298		0.02728	2.600	0.0101	0.198
14.00	1.26	0.84	0.132		0.08059		4.41E-04			0.02976	2.576	9.95E-03	0.194
14.40	1.25	0.84	0.132		0.08679	3.408	4.32E-04	0.298		0.03224	2.598	1.16E-02	0.197
14.80	1.25	0.84	0.133		0.09299		2.44E-04			0.03472	2.607	1.68E-02	0.199
15.00	1.25	0.84	0.133		0.09919	3.409	3.23E-04	0.299		0.03720	2.613	1.54E-02	0.199
15.60	1.25	0.85	0.134		0.1116	3.409	2.87E-04	0.299		0.04463		1.17E-02	
16.00	1.27	0.85	0.134		0.1240	3.410	2.71E-04	0.299		0.04959	2.627	3.58E-03	0.201
16.50	1.28	0.89	0.145		0.2480	3.417	2.67E-04	0.299		0.05703		8.65E-04	
					0.3720	3.427	1.90E-04	0.301					

Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)	Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)	Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)
5.00	2.58	2.20	0.416	23.00	0.73	0.24	0.043	0.7	7.00	0.24	0.563
5.20	2.52	2.44	0.450	23.60	0.80	0.26	0.033	0.8	7.23	0.48	0.574
5.40	2.31	2.61	0.480	24.00	0.80	0.26	0.034	0.9	7.48	0.94	0.589
5.60	2.06	2.67	0.501	24.60	0.82	0.25	0.029	1.0	7.70	1.56	0.606
5.80	1.83	2.63	0.510	25.00	0.83	0.25	0.026	1.2	6.99	2.22	0.593
6.00	1.63	2.56	0.515	25.60	0.86	0.24	0.022	1.4	7.11	2.46	0.604
6.20	1.48	2.45	0.512	26.00	0.88	0.25	0.022	1.6	6.75	2.91	0.606
6.40	1.37	2.33	0.504	26.60	0.87	0.26	0.023	1.8	6.89	3.70	0.637
6.60	1.29	2.22	0.492	27.00	0.87	0.25	0.022	2.0	4.67	4.67	0.654
6.80	1.23	2.11	0.478	27.60	0.89	0.23	0.019	2.2	4.94	5.16	0.681
7.00	1.18	2.01	0.462	28.00	0.90	0.23	0.017	2.4	3.94	5.08	0.686
7.20	1.15	1.91	0.445	28.60	0.91	0.22	0.015	2.6	3.25	4.77	0.681
7.40	1.13	1.82	0.425	29.00	0.92	0.22	0.014	2.8	2.73	4.42	0.674
7.60	1.12	1.75	0.406	29.60	0.94	0.22	0.014	3.0	2.30	4.16	0.674
7.80	1.11	1.68	0.390	30.00	0.95	0.22	0.014	3.5	1.69	3.44	0.646
8.00	1.11	1.61	0.370	31.00	0.97	0.23	0.014	4.0	1.33	2.64	0.571
8.20	1.12	1.55	0.350	32.00	0.98	0.24	0.015	4.5	1.32	1.96	0.428
8.40	1.13	1.50	0.332	33.00	0.98	0.25	0.015	5.0	1.63	1.60	0.312
8.60	1.14	1.45	0.317	34.00	0.99	0.25	0.016	5.5	1.72	1.57	0.302
8.80	1.17	1.41	0.301	35.00	0.99	0.26	0.017	6.0	1.73	1.45	0.276
9.00	1.19	1.40	0.294	36.00	0.99	0.27	0.018	6.5	1.78	1.36	0.257
9.20	1.21	1.38	0.289	37.00	0.99	0.28	0.019	7.0	1.83	1.36	0.257
9.40	1.21	1.38	0.287	38.00	0.98	0.28	0.021	7.5	1.72	1.51	0.289
9.60	1.21	1.38	0.285	39.00	0.97	0.29	0.022	8.0	1.54	1.37	0.260
9.80	1.21	1.37	0.285	40.00	0.95	0.29	0.023	8.5	1.55	1.23	0.226
10.00	1.20	1.37	0.286					9.0	0.99	0.93	0.179
10.20	1.19	1.37	0.286	<i>Tellurium, $\vec{E} \parallel c^{25}$</i>				9.5	1.47	1.25	0.233
10.40	1.18	1.37	0.287	0.01364	4.82	0.118	0.431	10.0	0.86	0.86	0.181
10.60	1.16	1.36	0.288	0.01488	5.26	0.0505	0.463	11.0	0.80	0.77	0.165
10.80	1.15	1.36	0.289	0.01612	5.47	0.0278	0.477	12.0	0.79	0.76	0.164
11.00	1.13	1.35	0.290	0.01736	5.59	0.0174	0.485	14.0	0.67	0.59	0.146
11.20	1.11	1.35	0.292	0.01860		0.0796		16.0	0.59	0.49	0.147
11.40	1.09	1.34	0.293	0.01984		0.0696		18.0	0.48	0.31	0.160
11.60	1.07	1.33	0.294	0.02108		0.0749		20.0	0.74	0.20	0.035
11.80	1.05	1.32	0.295	0.02232		0.1900		22.0	0.83	0.18	0.018
12.00	1.02	1.31	0.296	0.02356		0.2220		24.0	0.85	0.15	0.013
12.20	1.00	1.29	0.295	0.02480		0.0716		26.0	0.87	0.12	0.009
12.40	0.98	1.28	0.294	0.02604		0.0682		28.0	0.89	0.090	0.006
12.60	0.96	1.26	0.292	0.02728		0.0832		30.0	0.90	0.045	0.003
12.80	0.94	1.24	0.289	0.02976		0.0149					
13.00	0.93	1.22	0.286	0.03224		2.14E-03		<i>Tellurium, $\vec{E} \perp c^{25}$</i>			
13.60	0.91	1.16	0.272	0.03472		1.71E-02		0.01364	2.61	0.2980	0.204
14.00	0.90	1.15	0.272	0.03720	5.94	3.71E-03	0.507	0.01488	3.65	0.0894	0.325
14.60	0.85	1.15	0.285	0.03968		2.44E-03		0.01612	4.10	0.0535	0.370
15.00	0.80	1.13	0.293	0.04339	5.96	1.59E-03	0.508	0.01736	4.63	0.4990	0.420
15.60	0.72	1.08	0.301	0.04711		7.85E-04		0.01860		0.1170	
16.00	0.68	1.04	0.304	0.05083		7.38E-04		0.01984		0.0343	
16.60	0.63	0.97	0.301	0.05579		3.89E-04		0.02108	(4.42)	0.0421	0.398
17.00	0.60	0.92	0.296	0.06199	5.98	3.09E-04	0.509	0.02232		0.1060	
17.60	0.60	0.92	0.296	0.07439		2.52E-04		0.02356		0.0880	
18.00	0.55	0.79	0.274	0.08679		2.96E-04		0.02480		0.0458	
18.60	0.53	0.71	0.254	0.09919		3.68E-04		0.02604		0.0928	
19.00	0.53	0.65	0.236	0.12400	6.246	3.34E-04	0.524	0.02728		0.0886	
19.60	0.53	0.57	0.207	0.15500	6.253		0.525	0.02976		0.0232	
20.00	0.54	0.52	0.185	0.20660	6.286		0.526	0.03224		3.06E-03	
20.60	0.55	0.44	0.153	0.24800	6.316	7.48E-05	0.528	0.03472		1.25E-02	
21.00	0.57	0.39	0.127	0.31	6.372	1.18E-05	0.531	0.03720	4.71	2.65E-03	0.422
21.60	0.64	0.34	0.089	0.35		4.93E-04		0.03968		1.89E-03	
22.00	0.64	0.32	0.081	0.41		6.74E-03		0.04339	4.74	1.41E-03	0.425
22.60	0.69	0.27	0.058	0.5	6.53	2.30E-02	0.539	0.04711		8.38E-04	
				0.6	6.71	7.50E-02	0.549	0.05083		6.79E-04	

Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)	Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)	Energy (eV)	<i>n</i>	<i>k</i>	<i>R</i> ($\phi = 0$)
<i>Tungsten</i> ²⁷				5.40	2.92	3.58	0.586	22.80	0.49	0.69	0.272
0.10	14.06	54.71	0.983	5.60	2.43	3.70	0.618	23.20	0.49	0.66	0.263
0.20	3.87	28.30	0.981	5.80	2.00	3.61	0.637	23.60	0.48	0.62	0.252
0.25	2.56	22.44	0.980	6.00	1.70	3.42	0.643	24.00	0.49	0.57	0.234
0.30	1.83	18.32	0.979	6.20	1.47	3.24	0.646	24.40	0.50	0.53	0.213
0.34	1.71	15.71	0.973	6.40	1.32	3.04	0.640	24.80	0.51	0.49	0.191
0.38	1.86	13.88	0.963	6.60	1.21	2.87	0.631	25.20	0.53	0.46	0.171
0.42	1.92	12.63	0.954	6.80	1.12	2.70	0.619	25.60	0.55	0.43	0.150
0.46	1.69	11.59	0.952	7.00	1.06	2.56	0.607	26.00	0.57	0.40	0.132
0.50	1.40	10.52	0.952	7.20	1.01	2.43	0.593	26.40	0.59	0.38	0.117
0.54	1.23	9.45	0.948	7.40	0.98	2.30	0.573	26.80	0.61	0.37	0.105
0.58	1.17	8.44	0.938	7.60	0.95	2.18	0.556	27.00	0.62	0.36	0.099
0.62	1.28	7.52	0.917	7.80	0.93	2.06	0.533	27.50	0.64	0.34	0.085
0.66	1.45	6.78	0.888	8.00	0.94	1.95	0.505	28.00	0.67	0.32	0.073
0.70	1.59	6.13	0.856	8.20	0.94	1.86	0.481	28.50	0.69	0.31	0.065
0.74	1.83	5.52	0.810	8.40	0.96	1.76	0.449	29.00	0.71	0.30	0.057
0.78	2.12	5.00	0.759	8.60	0.99	1.70	0.422	29.50	0.73	0.30	0.052
0.82	2.36	4.61	0.710	8.80	1.01	1.65	0.401	30.00	0.75	0.29	0.047
0.86	2.92	4.37	0.661	9.00	1.01	1.60	0.388	31.00	0.78	0.29	0.042
0.90	3.11	4.44	0.660	9.20	1.02	1.55	0.369	32.00	0.79	0.29	0.040
0.94	3.15	4.43	0.658	9.40	1.03	1.50	0.352	33.00	0.82	0.28	0.033
0.98	3.15	4.36	0.653	9.60	1.05	1.44	0.329	34.00	0.84	0.29	0.032
1.00	3.14	4.32	0.649	9.80	1.09	1.38	0.307	35.00	0.85	0.31	0.033
1.10	3.05	4.04	0.627	10.00	1.13	1.34	0.287	36.00	0.85	0.32	0.036
1.20	3.00	3.64	0.590	10.20	1.19	1.33	0.274	37.00	0.84	0.33	0.039
1.30	3.12	3.24	0.545	10.40	1.24	1.34	0.270	38.00	0.83	0.33	0.040
1.40	3.29	2.96	0.515	10.60	1.27	1.36	0.274	39.00	0.81	0.33	0.042
1.50	3.48	2.79	0.500	10.80	1.29	1.39	0.282	40.00	0.80	0.33	0.045
1.60	3.67	2.68	0.494	11.00	1.28	1.42	0.290	<i>Vanadium</i> ⁹			
1.70	3.84	2.79	0.507	11.20	1.27	1.44	0.297	0.10	12.83	45.89	0.978
1.80	3.82	2.91	0.518	11.40	1.25	1.46	0.305	0.20	3.90	24.30	0.975
1.90	3.70	2.94	0.518	11.60	1.22	1.48	0.313	0.28	2.13	17.35	0.973
2.00	3.60	2.89	0.512	11.80	1.20	1.48	0.318	0.36	1.54	13.32	0.966
2.10	3.54	2.84	0.506	12.00	1.16	1.48	0.323	0.44	1.28	10.74	0.957
2.20	3.49	2.76	0.497	12.40	1.10	1.47	0.329	0.52	1.16	8.93	0.945
2.30	3.49	2.72	0.494	12.80	1.04	1.44	0.333	0.60	1.10	7.59	0.929
2.40	3.45	2.72	0.493	13.20	0.98	1.40	0.332	0.68	1.07	6.54	0.909
2.50	3.38	2.68	0.487	13.60	0.94	1.35	0.325	0.76	1.08	5.67	0.882
2.60	3.34	2.62	0.480	14.00	0.91	1.28	0.312	0.80	1.10	5.30	0.864
2.70	3.31	2.55	0.472	14.40	0.90	1.23	0.296	0.90	1.18	4.50	0.811
2.80	3.31	2.49	0.466	14.80	0.90	1.17	0.276	1.00	1.34	3.80	0.730
2.90	3.32	2.45	0.461	15.20	0.93	1.13	0.255	1.10	1.60	3.26	0.632
3.00	3.35	2.42	0.459	15.60	0.97	1.12	0.246	1.20	1.93	2.88	0.543
3.10	3.39	2.41	0.460	16.00	0.98	1.14	0.249	1.30	2.25	2.71	0.498
3.20	3.43	2.45	0.465	16.40	0.97	1.17	0.260	1.40	2.48	2.72	0.491
3.30	3.45	2.55	0.476	16.80	0.94	1.19	0.273	1.50	2.57	2.79	0.499
3.40	3.39	2.66	0.485	17.20	0.90	1.21	0.289	1.60	2.57	2.84	0.507
3.50	3.24	2.70	0.488	17.60	0.85	1.21	0.304	1.70	2.52	2.88	0.512
3.60	3.13	2.67	0.482	18.00	0.80	1.20	0.317	1.80	2.45	2.88	0.515
3.70	3.05	2.62	0.476	18.40	0.74	1.18	0.330	1.90	2.36	2.85	0.514
3.80	2.99	2.56	0.468	18.80	0.69	1.15	0.340	2.00	2.34	2.81	0.509
3.90	2.96	2.50	0.460	19.20	0.64	1.11	0.347	2.10	2.31	2.78	0.506
4.00	2.95	2.43	0.451	19.60	0.60	1.07	0.353	2.20	2.28	2.80	0.510
4.20	3.02	2.33	0.440	20.00	0.56	1.02	0.354	2.30	2.23	2.83	0.516
4.40	3.13	2.32	0.442	20.40	0.54	0.97	0.350	2.40	2.15	2.88	0.528
4.60	3.24	2.41	0.455	20.80	0.52	0.92	0.342	2.50	2.02	2.91	0.540
4.80	3.33	2.57	0.475	21.20	0.50	0.87	0.331	2.60	1.89	2.92	0.552
5.00	3.40	2.85	0.505	21.60	0.50	0.82	0.318	2.70	1.74	2.89	0.561
5.20	3.27	3.27	0.548	22.00	0.49	0.77	0.303	2.80	1.61	2.85	0.569
				22.40	0.49	0.73	0.287	2.90	1.48	2.80	0.577

Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$	Energy (eV)	n	k	$R(\phi = 0)$
3.60	1.10	0.74	0.013	9.00	1.65	0.91	0.025	17.20	1.09	0.74	0.013
3.70	1.07	0.73	0.013	9.20	1.63	0.90	0.025	17.60	1.13	0.75	0.013
3.80	1.04	0.72	0.012	9.40	1.60	0.89	0.024	18.00	1.17	0.76	0.014
3.90	1.01	0.71	0.012	9.60	1.57	0.89	0.023	18.40	1.21	0.78	0.014
4.00	0.98	0.70	0.012	9.80	1.52	0.87	0.021	18.80	1.24	0.79	0.014
4.20	0.94	0.68	0.013	10.00	1.47	0.86	0.020	19.20	1.27	0.80	0.015
4.40	0.89	0.67	0.013	10.20	1.42	0.84	0.018	19.60	1.29	0.80	0.015
4.60	0.85	0.65	0.014	10.40	1.35	0.82	0.016	20.00	1.30	0.81	0.015
4.80	0.81	0.64	0.014	10.50	1.32	0.81	0.016	20.60	1.29	0.80	0.015
5.00	0.78	0.63	0.015	10.60	1.28	0.80	0.015	21.00	1.27	0.80	0.015
5.20	0.77	0.62	0.016	10.80	1.23	0.78	0.014	21.60	1.23	0.78	0.014
5.40	0.77	0.62	0.016	11.00	1.19	0.77	0.014	22.00	1.20	0.77	0.014
5.60	0.80	0.63	0.014	11.20	1.16	0.76	0.013	22.60	1.15	0.76	0.013
5.80	0.87	0.66	0.013	11.40	1.13	0.75	0.013	23.00	1.12	0.75	0.013
6.00	1.00	0.71	0.012	11.60	1.11	0.74	0.013	23.60	1.08	0.73	0.013
6.20	1.11	0.75	0.013	11.80	1.09	0.74	0.013	24.00	1.05	0.73	0.013
6.40	1.23	0.78	0.014	12.00	1.08	0.73	0.013	24.60	1.02	0.71	0.012
6.60	1.33	0.81	0.016	12.40	1.05	0.72	0.012	25.00	1.00	0.71	0.012
6.80	1.42	0.84	0.018	12.80	1.01	0.71	0.012	25.60	0.97	0.69	0.012
7.00	1.49	0.86	0.020	13.20	0.98	0.70	0.012	26.00	0.95	0.69	0.013
7.20	1.54	0.88	0.022	13.60	0.95	0.69	0.013	26.60	0.91	0.67	0.013
7.40	1.58	0.89	0.023	14.00	0.92	0.68	0.013	27.00	0.88	0.66	0.013
7.60	1.61	0.90	0.024	14.40	0.89	0.67	0.013	27.60	0.84	0.65	0.014
7.80	1.63	0.90	0.025	14.80	0.90	0.67	0.013	28.00	0.83	0.64	0.014
8.00	1.66	0.91	0.026	15.20	0.92	0.68	0.013	28.60	0.82	0.64	0.014
8.20	0.67	0.91	0.026	15.60	0.95	0.69	0.013	29.00	0.81	0.64	0.014
8.40	1.68	0.92	0.026	16.00	0.98	0.70	0.012	29.60	0.82	0.64	0.014
8.60	1.68	0.92	0.026	16.40	1.01	0.71	0.012	30.00	0.82	0.64	0.014
8.80	1.66	0.91	0.026	16.80	1.04	0.72	0.012				

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