

# Strontium Fluoride (SrF<sub>2</sub>)

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Strontium fluoride offers a wide range of transparency, from the ultraviolet to the long-wave infrared (0.13 to 11  $\mu\text{m}$ ), with low reflectance loss and low dispersion. This combination of broad transparency and low dispersion is rare in optical-window materials used in infrared systems. Like BaF<sub>2</sub>, it allows experimenters to conveniently perform initial system alignment in the visible and then switch to the IR with only minor correction of the optical components. Furthermore, strontium fluoride has good surface hardness (Knoop scale 150 kgf/cm<sup>2</sup>) and is not hygroscopic. It has a fracture strength between those of CaF<sub>2</sub> (90 MPa) and BaF<sub>2</sub> (27 MPa).

Strontium fluoride is a cubic crystal with the fluorite structure, space group Fm $\bar{3}m$  ( $O_h^h$ ), with four formula units per unit cell. The lattice constant is 5.7996 Å, giving a theoretical density of 4.277 g/cm<sup>3</sup>. The atomic arrangement in strontium fluoride occupy the 4(a) sites (m $\bar{3}m$  or  $O_h$  symmetry), and the fluoride ions occupy the 8(c) sites ( $\bar{4}3m$  or  $T_d$  symmetry). Melting point is 1750 K.

Ultraviolet reflectance of bulk SrF<sub>2</sub> was measured by Genie et al. [1] up to 20 eV at room temperature and liquid-helium temperature. Shuboff [2] measured the room-temperature reflectance up to 50 eV. Unfortunately, these measurements are not indexed to optical constants. However, similar measurements from 10 to 50 eV by Nisar and Robin [3] later were reduced to optical constants using the Kramers-Kronig relation. These data are listed in Table I and plotted in Fig. 1.

The Urbach tail represents absorption below the band gap, defining the end of transparency. The room-temperature absorption coefficient was determined from transmission measurements by Towhid and Mlynek [4] in the spectral range from 9.3 to 9.8 eV and fitted to the functional form

$$\beta_{\text{Urb}}(E, T) = \beta_{\text{Urb}} \exp(\alpha_{\text{Urb}}(T)(E - E_g)/k_B T), \quad (1)$$

where  $\beta_{\text{Urb}}$  is the absorption coefficient (typically in units of cm<sup>-1</sup>),  $\beta_{\text{Urb}}$  is

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a scaling coefficient (in units of  $\text{cm}^{-1}$ ),  $E_p$  is the band-gap energy at absolute zero temperature (typically given in units of eV),  $k_B$  is Boltzmann's constant, and  $T$  is temperature in Kelvin. The exponential factor  $\alpha_s(T)$  is given by the equation

$$\alpha_s(T) = \alpha_0 \frac{2k_B T}{E_p} \tanh \frac{E_p}{2k_B T} \quad (2)$$

where  $E_p$  is an effective acoustic-phonon energy of the material. The Urbach tail parameters for  $\text{SrF}_2$  are  $\beta_{\text{UD}} \approx 1.35 \times 10^8 \text{ cm}^{-1}$ ,  $E_p = 10.670 \text{ eV}$ ,  $\alpha_0 = 0.60$ , and  $E_g = 0.044 \text{ eV}$ . The room-temperature (295 K) values for  $k$  in Table I and Fig. 1 for this spectral region are generated using Eqs. (1) and (2).

The room-temperature (25°C) index of refraction from the ultraviolet to the infrared (0.22 to 11.5  $\mu\text{m}$ ) has been measured by Feldman *et al.* [5] using a prism and the minimum-deviation technique. A Sellmeier model of the form

$$n^2(\lambda, T) = 1 + \sum_j \frac{A^2 d\epsilon_j(T)}{\lambda^2 - \lambda_j^2(T)} \quad (3)$$

is used to represent the data accurately. Model parameters are listed in Table II. A comprehensive analysis on a variety of data sets was performed by Li [6]. A Sellmeier formula was also generated covering the range from 0.15 to 14  $\mu\text{m}$  at  $T = 20^\circ \text{C}$ . The Li model parameters are also listed in Table II. Because the Sellmeier model has a physical basis, the extrapolation to higher frequencies beyond the experimental limit is justified. The values for  $n$  in Table I and Fig. 1 for the spectral regions indicated are generated by the appropriate formula. A temperature-dependent Sellmeier model covering the range from 100 to 450  $\mu\text{m}$  has been developed by Tropf [7]. Near-room-temperature thermo-optic coefficients measured at different wavelengths are listed in Table III [5, 8-10]. Notice that the thermo-optic coefficients are negative. This gives  $\text{SrF}_2$  some unusual performance as an optical element. An unusual material has the property that

$$\frac{d(n(T)L(T))}{dT} = 0.$$

The value of the optical path derivative as given earlier for  $\text{SrF}_2$  is  $6.9 \times 10^{-6} \text{ eL per Kelvin}$  at visible and near infrared wavelengths. Thus,  $\text{SrF}_2$  is useful for optical designs that require insensitivity to temperature.

Low-absorptivity data [11] at DF (3.4  $\mu\text{m}$ ) and HP (3.7  $\mu\text{m}$ ) laser wavelengths indicate low-level absorption is roughly the width of the transparency range of  $\text{SrF}_2$ . Material obtained from Optosys had absorption-

coefficient values of  $2.0 \times 10^{-4} \text{ cm}^{-1}$  at the DF laser wavelength and  $1.4 \times 10^{-4} \text{ cm}^{-1}$  at the EF laser wavelength. The corresponding  $k$  values are listed in Table I and plotted in Fig. 1. The low-level absorption and the near athermal performance of SrF<sub>2</sub> make it a candidate window material for high-power laser applications.

The multiphonon absorption (multiple-quantum lattice vibrations) edge marks the end of infrared transparency. Absorption-coefficient measurements by Demich [12] was represented by the simple formula

$$\beta_{\text{MFA}}(\nu) = \beta_0 \exp\left(-\gamma \frac{\nu}{\nu_0}\right), \quad (6)$$

where for SrF<sub>2</sub>,  $\beta_0 = 22,548 \text{ cm}^{-1}$  and  $\nu_0/\gamma = 90.4 \text{ cm}^{-1}$  at room temperature. The values of  $k$  in the spectral range from 830 to 1300  $\text{cm}^{-1}$  in Table I and Fig. 1 are obtained from this formula. The data cover the three-phonon to four-phonon regions. Temperature-dependent experimental data on the absorption coefficient is described in Lipson *et al.* [13]. A temperature-dependent multiphonon model has been developed by Thomas *et al.* [14] and applied to SrF<sub>2</sub>. Measurements in the two-phonon region was reported by Kaber *et al.* [15]. The experimental results are listed in Table I and plotted in Fig. 1.

The one-phonon (one-quantum three-dimensional lattice vibrations) regime is opaque and therefore characterized by reflectance measurements. A room-temperature reflectance measurement is reported by Kaim *et al.* [15]. The classical-oscillator model is often used to fit the reflectance data and then to derive the optical constants. The classical-oscillator model is expressed in terms of the relative permittivity,  $\epsilon_r(\nu, T)$ , as given by<sup>1</sup>

$$\epsilon_r(\nu, T) = \epsilon_\infty(T) + \sum_j \frac{\Delta\epsilon_j(T) \nu_j^2(T)}{\nu_j^2(T) - \nu^2 + j\Gamma_j(\nu, T)\nu}, \quad (5)$$

where  $\Delta\epsilon_j(T)$ ,  $\Gamma_j(\nu, T)$  and  $\nu_j(T)$  are the  $j$ th-mode strength, linewidth, and long-wavelength omnivector optical frequency,  $\nu_{TO}$ , respectively. The sum on  $j$  is over all transparent optical modes. For SrF<sub>2</sub>, there is only one infrared-active vibrational mode based on group theory of a perfect lattice. This is formally stated as

$$\Gamma = F_{1g}(\text{IR}) + F_{2g}(\text{R}).$$

The first mode is infrared active (mode 1 in Table IV), and the second mode is Raman active (286  $\text{cm}^{-1}$  [16]). Table IV lists the classical-oscillator model parameters used to fit the experimental data. More than one mode is needed

<sup>1</sup> The equation was the engineering convention for a time-harmonic field of  $\exp(-i\omega t)$ . To convert to the convention commonly used in the IEC units, we let  $i = -j$ .

to account for impurities and defects in real materials. Also, the high-frequency edge of the reflectance spectrum must include two-phonon contributions (see mode 2 in Table IV). Unfortunately, the value of  $\epsilon_\infty$  used by Kasper *et al.* is not consistent with the value from the Sellmeier models mentioned previously, and the same is true for the static dielectric constant. For this reason the model parameters have been modified to agree with these limiting values and still match the reflectance spectrum. These modified values are listed in Table IV. The  $n$  and  $k$  values in Table I and Fig. 1 are generated by Eq. (5) and the modified parameters in Table IV.

Below  $\lambda_{\text{TO}}$ , the index of refraction,  $n$ , is determined by extrapolation of the classical-oscillator model with reasonable accuracy. The absorption coefficient has been measured by Boucsworth (17). Table I and Fig. 1 list and display the resulting values. The values are above the one-phonon red wing because multiphonon difference bands are also included in this spectral region.

The static-dielectric-constant data is the sum of all the strengths of higher-frequency oscillators including vibrational and electronic. Using Eq. (5), the static dielectric constant,  $\epsilon_s(T) = \epsilon'(0, T)$ , becomes

$$\epsilon_s(T) = \epsilon_\infty(T) + \sum_j \Delta\epsilon_j(T). \quad (6)$$

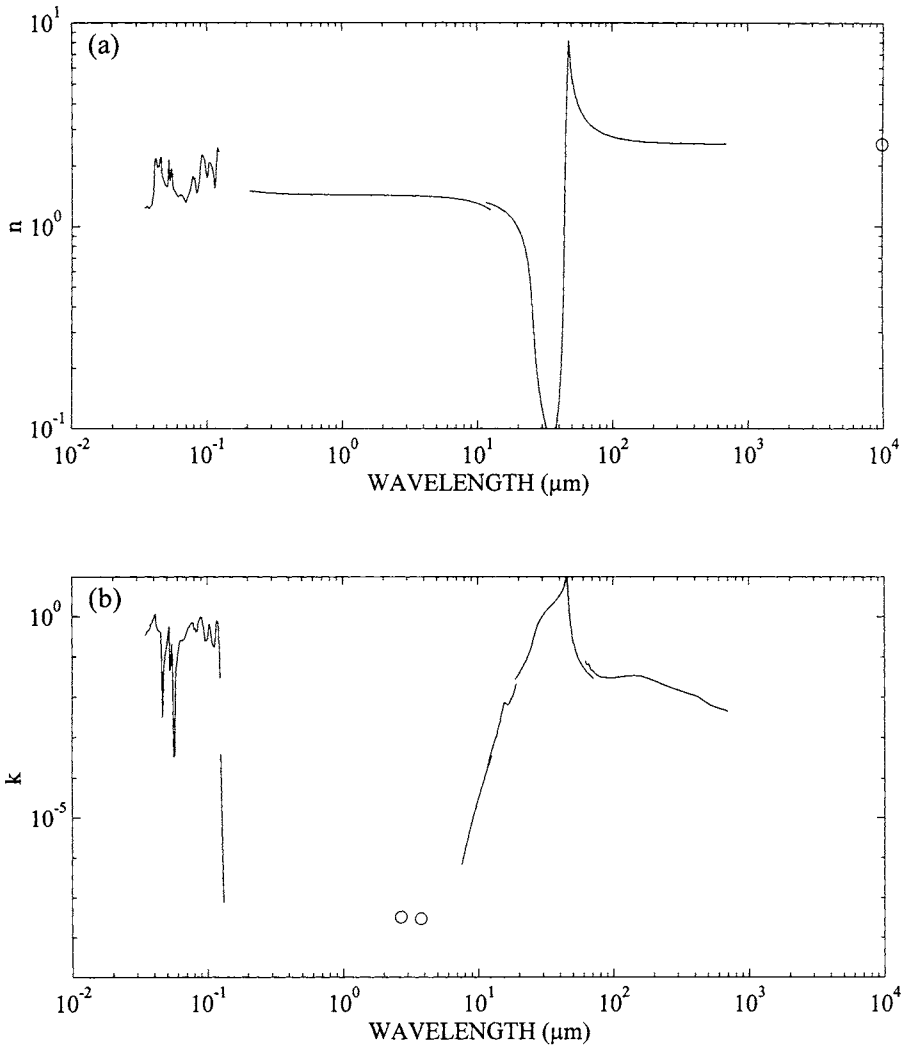
Based on the parameters listed in Table IV, we expect the value of  $\epsilon_s$  to be between 6.14 and 6.5. Experimental results, obtained from capacitance bridges operating from 1 kHz to 1 MHz, produce values from 6.45 to 6.48, with most results closer to the higher value [18-21]. The corresponding refractive index value is  $n_s = 2.543$ . The static-index values inaccurately are located at  $1 \times 10^4 \mu\text{m}$ ; however, there is little difference in the value of  $n_s$  at the measured wavelength and the plotted wavelength.

Whenever possible, physically based models are used to improve experimental measurements in Table I and Fig. 1. The models tend to reduce noise and allow interpolation and extrapolation to obtain meaningful results where no experimental data exists. In some cases, the models discussed in this section also allow temperature-dependent data to be obtained.

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**Fig. 1.** (a) Log-log plot of the room-temperature index of refraction,  $n$ , for  $\text{SrF}_2$  versus wavelength in microns. (b) Log-log plot of the room-temperature extinction coefficient,  $k$ , for  $\text{SrF}_2$  versus wavelength in microns. Isolated data points are indicated with a symbol.

TABLE I  
 Values of  $\alpha$  and  $k$  for Boronhan Fluoride from Various References\*

$\nu$	$\text{cm}^{-1}$	$\mu\text{m}$	$\alpha$	$k$
25.05	229500	0.00454	1.242	0.3443 [3]
25.24	264300	0.00378	1.236	0.3963
26.04	367800	0.00272	1.241	0.4173
26.38	376900	0.00265	1.260	0.4271
26.83	371800	0.00269	1.236	0.4096
28.15	357300	0.00280	1.225	0.3556
28.27	359000	0.00279	1.237	0.3975
28.77	353900	0.00283	1.261	0.4290
29.12	344000	0.00291	1.272	0.4971
29.34	353900	0.00283	1.278	0.7205
31.84	296000	0.00338	1.293	0.7973
31.21	321700	0.00311	1.407	0.7007
31.03	290800	0.00344	1.448	0.9406
30.50	260800	0.00383	1.460	1.011
30.74	166500	0.00601	1.578	1.064
30.58	262400	0.00381	1.064	1.115
30.34	262700	0.00381	1.931	1.137
30.33	264700	0.00378	2.051	1.112
30.30	266800	0.00375	2.025	1.063
30.26	264000	0.00382	2.109	1.467
30.21	262800	0.00384	2.137	0.9336
30.13	262800	0.00384	2.154	0.4243
30.06	262500	0.00385	2.146	0.3007
29.97	261700	0.00387	2.145	0.7403
29.86	260800	0.00389	2.100	0.6764
29.76	260000	0.00391	2.196	0.6386
29.64	259000	0.00394	2.140	0.5702
28.80	232700	0.00429	1.931	0.4333
28.58	240300	0.00416	1.974	0.4471
28.35	226700	0.00441	1.924	0.4138
28.18	227000	0.00440	1.969	0.4252
27.68	223300	0.00448	2.180	0.6053
27.43	231200	0.00432	2.195	0.1572
27.33	220300	0.00454	2.077	0.0794
27.08	218200	0.00463	2.075	0.00111
26.37	212000	0.00472	1.279	0.00483
26.17	213900	0.00468	1.740	0.05753
26.05	210800	0.00474	1.707	0.00398
25.56	203700	0.00491	1.604	0.1553
25.51	201700	0.00496	1.699	0.1377
24.50	199100	0.00502	1.398	0.2805
24.42	196800	0.00509	1.946	0.7489
24.36	195200	0.00512	1.613	0.4236
24.30	193400	0.00517	1.799	0.2621

(continued)

\* References given in brackets.

TABLE I (Continued)  
Stochastic Fluctuosity

$\nu$	$\text{cm}^{-1}$	$\mu\text{m}$	$n$	$f$
23.99	193900	0.005169	2.043	0.5311
23.99	193000	0.005180	2.110	0.4473
23.77	191700	0.005237	2.129	0.4940
23.59	189500	0.005278	2.746	0.8900
23.39	187100	0.005344	1.580	0.5006
23.22	184900	0.005389	1.820	0.2049
22.73	180300	0.005449	1.910	0.1022
22.64	180000	0.005476	1.925	0.0627
22.44	181000	0.005528	1.813	0.0253
22.13	178300	0.005403	1.292	0.0063
21.77	175000	0.005589	1.923	0.0087
21.36	170700	0.007730	1.980	0.0216
21.29	173700	0.005523	1.471	0.0193
20.98	169000	0.005916	1.447	0.0027
20.20	162000	0.008133	1.407	0.0047
19.73	159000	0.008291	1.434	0.2916
19.45	157000	0.008370	1.431	0.2237
19.17	154000	0.008408	1.431	0.2530
18.26	149100	0.008973	1.419	0.2518
18.45	148400	0.008713	1.367	0.2730
18.21	149500	0.008408	1.564	0.3070
22.78	145500	0.008863	1.319	0.3907
17.39	161800	0.007408	1.319	0.4436
22.40	140400	0.007114	1.369	0.4883
17.28	170300	0.007181	1.383	0.3240
17.00	159400	0.007368	1.411	0.3437
16.83	159400	0.007368	1.437	0.3801
16.34	153800	0.007380	1.354	0.5087
16.13	150700	0.007880	1.200	0.4828
15.92	148300	0.007790	1.708	0.6637
15.76	157200	0.007965	1.709	0.5848
15.71	155700	0.007882	1.742	0.4449
15.69	155900	0.007905	1.760	0.5019
16.64	138700	0.007020	1.742	0.4731
16.43	134400	0.008042	1.708	0.5724
16.29	128000	0.008159	1.737	0.4817
15.13	122200	0.008286	1.746	0.4011
14.30	159000	0.008578	1.591	0.4348
14.43	115000	0.008472	1.432	0.3739
14.37	117800	0.008308	1.461	0.0043
14.20	120000	0.008423	1.520	0.3740
14.27	115400	0.008283	1.584	0.3947
14.20	114000	0.007736	1.601	0.6253
14.07	119800	0.008614	1.671	0.9229
14.00	118900	0.008408	1.736	0.9630
14.74	110800	0.008023	1.693	0.6530



TABLE I (Continued)  
 Strontium Fluoride

$\nu$	$\text{cm}^{-1}$	$\rho_{\text{sr}}$	$n$	$k$
13.60	109709	0.09116	2.185	0.2993
13.53	109309	0.09180	2.225	0.2977
13.43	108939	0.09251	2.249	0.2974
13.38	107909	0.09298	2.294	0.2972
13.31	107309	0.09317	2.341	0.2983
13.11	105909	0.09434	2.384	0.2948
13.04	105509	0.09452	2.185	0.2994
12.94	99909	0.1009	1.738	0.2938
12.28	99709	0.1010	1.779	0.3342
12.24	93759	0.1013	1.720	0.2948
12.19	92569	0.1017	1.766	0.2998
12.14	92939	0.1021	1.809	0.3389
12.04	92549	0.1024	1.818	0.4044
12.02	92299	0.1032	1.991	0.4913
11.92	96129	0.1083	2.059	0.3266
11.87	95709	0.1085	2.080	0.3012
11.81	92389	0.1090	2.080	0.4922
11.71	94669	0.1099	2.043	0.3738
11.58	93399	0.1091	2.029	0.3915
11.48	92529	0.1093	2.035	0.3326
11.52	91279	0.1096	1.993	0.3983
11.49	89409	0.1108	1.987	0.4687
10.91	88819	0.1138	1.998	0.2072
10.84	87409	0.1143	1.999	0.2621
10.79	87009	0.1149	1.989	0.3263
10.76	86779	0.1138	1.984	0.3096
10.72	85449	0.1137	1.989	0.4378
10.67	86999	0.1182	1.972	0.4911
10.64	86349	0.1186	1.971	0.4739
10.58	85539	0.1172	1.789	0.2407
10.59	86989	0.1178	1.925	0.4912
10.58	85969	0.1206	2.415	0.4706
10.56	93009	0.1789	3.623	0.6086
10.22	82969	0.1207	2.449	0.2499
10.25	83699	0.1210	2.449	0.4914
10.24	82379	0.1271	2.611	0.4282
10.21	82969	0.1214	2.424	0.3630
10.17	83699	0.1219	2.602	0.2947
10.17	80789	0.1211	2.394	0.2826
10.10	81789	0.1239	2.379	0.1654
10.06	80099	0.1234	2.339	0.1017
10.01	80739	0.1238	2.315	0.0293
9.239	79989	0.1233	1.746 [6]	0.000305 [6]
9.091	79089	0.1269	1.720	0.2689 5
9.224	78339	0.1233	1.225	2.0288 5
9.049	79829	0.1238	1.716	1.8298 6

(continued)

TABLE I (Continued)  
Strontium Fluoride

eV	cm <sup>-1</sup>	μm	n	k
9.574	77220	0.1295	1.707	1.174E-6
9.501	76630	0.1305	1.699	2.922E-7
9.429	76050	0.1315	1.691	7.428E-8
8.266	66670	0.150	1.594	
7.085	57140	0.175	1.534	
6.199	50000	0.200	1.504	
5.510	44440	0.225	1.486	
5.904	47619	0.21	1.49513 [5]	
4.959	40000	0.25	1.47336	
3.999	32258	0.31	1.45725	
3.542	28571	0.35	1.45131	
3.024	24390	0.41	1.44556	
2.755	22222	0.45	1.44309	
2.431	19608	0.51	1.44029	
2.254	18182	0.55	1.43906	
2.032	16393	0.61	1.43740	
1.907	15384	0.65	1.43675	
1.746	14084	0.71	1.43560	
1.653	13333	0.75	1.43527	
1.531	12346	0.81	1.43435	
1.362	10989	0.91	1.43343	
1.228	9901.0	1.01	1.43269	
1.117	9009.0	1.11	1.43206	
1.025	8264.5	1.21	1.43151	
0.9464	7633.6	1.31	1.43100	
0.8793	7092.2	1.41	1.43051	
0.8211	6622.5	1.51	1.43003	
0.7701	6211.2	1.61	1.42956	
0.7251	5847.9	1.71	1.42909	
0.6850	5524.9	1.81	1.42861	
0.6491	5235.6	1.91	1.42812	
0.6168	4975.1	2.01	1.42761	
0.5876	4739.3	2.11	1.42709	
0.5610	4524.9	2.21	1.42656	
0.5367	4329.0	2.31	1.42601	
0.5145	4149.4	2.41	1.42544	
0.4940	3984.1	2.51	1.42485	
0.4750	3831.4	2.61	1.42424	
0.4590	3703.7	2.70		3.20E-8 [11]
0.4575	3690.0	2.71	1.42361	
0.4412	3558.7	2.81	1.42296	
0.4261	3436.4	2.91	1.42228	
0.4119	3322.3	3.01	1.42159	
0.3987	3215.4	3.11	1.42087	
0.3862	3115.3	3.21	1.42013	
0.3746	3021.1	3.31	1.41937	

TABLE I (Continued)

## Strontium Fluoride

eV	cm <sup>-1</sup>	μm	<i>n</i>	<i>k</i>
0.3636	2932.5	3.41	1.41858	
0.3532	2849.0	3.51	1.41778	
0.3434	2770.1	3.61	1.41694	
0.3342	2695.4	3.71	1.41609	
0.3260	2631.6	3.80		2.90E-8 [11]
0.3254	2624.7	3.81	1.41521	
0.3171	2557.5	3.91	1.41430	
0.3092	2493.8	4.01	1.41337	
0.3017	2433.1	4.11	1.41242	
0.2945	2375.3	4.21	1.41144	
0.2877	2320.2	4.31	1.41044	
0.2811	2267.6	4.41	1.40941	
0.2749	2217.3	4.51	1.40835	
0.2689	2169.2	4.61	1.40727	
0.2633	2123.1	4.71	1.40617	
0.2578	2079.0	4.81	1.40503	
0.2525	2036.7	4.91	1.40388	
0.2475	1996.0	5.01	1.40269	
0.2426	1956.9	5.11	1.40148	
0.2380	1919.4	5.21	1.40024	
0.2335	1883.2	5.31	1.39898	
0.2292	1848.4	5.41	1.39768	
0.2250	1814.9	5.51	1.39636	
0.2210	1782.5	5.61	1.39502	
0.2171	1751.3	5.71	1.39364	
0.2134	1721.2	5.81	1.39223	
0.2098	1692.1	5.91	1.39080	
0.2063	1663.9	6.01	1.38934	
0.2029	1636.7	6.11	1.38785	
0.1997	1610.3	6.21	1.38633	
0.1965	1584.8	6.31	1.38478	
0.1934	1560.1	6.41	1.38320	
0.1905	1536.1	6.51	1.38159	
0.1876	1512.9	6.61	1.37995	
0.1848	1490.3	6.71	1.37828	
0.1821	1468.4	6.81	1.37658	
0.1794	1447.2	6.91	1.37485	
0.1769	1426.5	7.01	1.37308	
0.1744	1406.5	7.11	1.37128	
0.1720	1387.0	7.21	1.36946	
0.1696	1368.0	7.31	1.36759	
0.1673	1349.5	7.41	1.36570	
0.1651	1331.6	7.51	1.36377	
0.1629	1314.1	7.61	1.36181	
0.1608	1297.0	7.71	1.35981	
0.1588	1280.4	7.81	1.35778	

(continued)

TABLE I (Continued)  
 Ground-State Parameters

$\nu$	$\text{cm}^{-1}$	$\mu$	$a$	$k$
0.1347	1264.2	7.91	1.31572	
0.1348	1268.4	8.01	1.31563	
0.1351	1254.6	8.1	1.3117 [8]	1.70335E-6 [12]
0.1440	1162.1	6.6	1.2400	4.11E-6
0.1543	1028.9	9.1	1.3283	3.348E-4
0.1289	1501.7	9.6	1.3151	1.71E-5
0.1229	909.1	10.1	1.3900	3.12E-3
0.1170	943.4	10.6	1.3854	3.41E-3
0.1117	929.9	11.1	1.3845	9.30E-5
0.1040	891.1	11.6	1.3701	0.0190
0.10610	857.7	11.83	1.3150	0.000173 [15]
0.1007	809.5	12.35	1.3503	0.00361
0.09746	787.7	12.45	1.3031	0.000302
0.09326	758.3	13.02	1.2825	0.000431
0.09346	723.0	13.27	1.3708	0.000792
0.09181	738.9	13.55	1.2825	0.000253
0.09557	723.2	[3.43]	1.3611	0.00120
0.09526	712.8	14.03	1.2353	0.01199
0.09025	684.4	15.07	1.2573	0.00232
0.09030	688.4	15.12	1.2805	0.00430
0.09344	646.7	15.81	1.2875	0.00719
0.09092	634.5	15.70	1.1979	0.00702
0.09796	628.7	15.91	1.1628	0.00705
0.09098	613.9	15.13	1.1640	0.00662
0.09098	611.7	15.58	1.1767	0.00642
0.09085	608.9	15.61	1.1659	0.00419
0.09280	587.2	17.03	1.1254	0.00603
0.06359	553.2	18.20	1.0991	0.0107
0.06440	528.1	19.12	1.0442	0.0908
0.06447	520.	18.23	1.094 <sup>b</sup>	0.0297 <sup>a</sup>
0.06323	510	19.61	1.015	0.0531
0.06199	500	20.09	0.991	0.0970
0.06015	480	20.41	0.970	0.0436
0.05991	480	20.89	0.934	0.0470
0.05927	470	21.35	0.911	0.0339
0.05792	460	21.74	0.884	0.0512
0.06050	480	22.22	0.822	0.0308
0.06440	460	22.78	0.774	0.0428
0.05391	430	23.26	0.719	0.0977
0.05207	420	23.81	0.673	0.111
0.05063	410	24.39	0.628	0.144
0.05025	400	25.00	0.586	0.189
0.04825	390	25.62	0.546	0.251
0.04711	380	26.33	0.508	0.378
0.04587	370	27.03	0.473	0.519
0.04465	360	27.79	0.441	0.683

<sup>a</sup> Results obtained from excited classical-oscillator parameters in Table IV.

TABLE I (Continued)

Bromine Fluoride

$\nu$	$\text{cm}^{-1}$	$\mu\text{m}$	$\tau$	$k$
0.04339	230	28.57	0.126	0.000
0.04215	240	29.11	0.137	0.040
0.04090	250	30.00	0.122	1.087
0.03966	260	31.25	0.125	1.243
0.03840	270	32.20	0.0998	1.413
0.03720	280	33.53	0.0925	1.687
0.03596	290	32.20	0.0881	1.818
0.03472	300	33.71	0.0674	2.071
0.03348	270	37.04	0.0923	2.370
0.03224	300	33.40	0.106	2.270
0.03100	290	40.00	0.130	3.394
0.02976	340	41.67	0.150	4.121
0.02852	290	43.88	0.448	3.894
0.02728	330	40.40	3.166	10.030
0.02606	210	47.62	4.152	1.013
0.02480	390	50.00	5.520	0.272
0.02356	180	52.63	4.534	0.136
0.02233	160	56.60	3.904	0.0825
0.02108	170	56.92	3.094	0.0504
0.01984	180	62.50	3.606	0.0445
0.01860	160	66.67	3.327	0.0440
0.01736	180	71.83	3.698	0.0282
0.01612	160.0	62.14	3.426	0.0302 [17]
0.01488	163.7	62.21	3.622	0.0766
0.01364	160.0	62.34	3.413	0.0706
0.01240	159.7	62.43	3.398	0.0725
0.01116	158.9	62.96	3.393	0.0667
0.00992	157.1	63.68	3.349	0.0643
0.00868	156.9	64.53	3.309	0.0620
0.00744	151.3	66.10	3.298	0.0629
0.00620	156.0	66.67	3.227	0.0543
0.00496	169.2	67.04	3.214	0.0290
0.00372	168.2	67.69	3.198	0.0529
0.00248	145.4	66.23	3.173	0.0472
0.00124	143.7	69.69	3.136	0.0454
0.00744	141.0	70.94	3.102	0.0634
0.00721	139.8	73.08	2.076	0.0610
0.00740	137.2	72.86	3.023	0.0383
0.01090	126.7	76.24	3.026	0.0362
0.01546	122.7	78.07	2.064	0.0329
0.01531	123.5	80.04	2.622	0.0315
0.01487	119.9	83.56	2.930	0.0304
0.01370	116.4	86.21	2.823	0.0295
0.01290	104.9	95.40	2.290	0.0204
0.01235	97.0	103.4	2.740	0.0229
0.01172	94.0	105.3	2.735	0.0204

(Continued)

TABLE I (Continued)  
Strontium Fluoride

eV	cm <sup>-1</sup>	μm	<i>n</i>	<i>k</i>
0.01102	88.9	112.5	2.710	0.0312
0.01044	84.2	118.8	2.690	0.0318
0.00978	78.9	126.8	2.670	0.0331
0.00924	74.5	134.2	2.655	0.0333
0.00872	70.3	142.2	2.642	0.0335
0.00824	66.5	150.4	2.630	0.0330
0.00769	62.0	161.3	2.618	0.0327
0.00720	58.1	172.2	2.608	0.0304
0.00664	53.6	186.6	2.598	0.0275
0.00595	48.0	208.4	2.587	0.0239
0.00534	43.1	232.1	2.578	0.0204
0.00481	38.8	258.0	2.571	0.0176
0.00359	29.0	344.8	2.559	0.0127
0.00293	23.7	422.7	2.554	0.0101
0.00239	19.3	519.0	2.550	0.0062
0.00178	14.4	697.0	2.547	0.0045
0.0	0.0	∞	2.543 [21]	0.0

TABLE II  
Room-Temperature Sellmeier Model Parameters

<i>i</i> th mode	$\Delta\epsilon_i$	$\lambda_i$
Feldman <i>et al.</i> model		
1	0.67805894	0.05628989
2	0.37140533	0.10801027
3	3.8484723	34.649040
Li model		
1	0.33973	0.0
2	0.70970	0.09597
3	0.1788	26.03
4	3.8796	45.60

**TABLE III**  
**Thermo-optical Coefficients of Strontium Fluoride**

Wavelength ( $\mu\text{m}$ )	$dn/dT$ ( $10^{-6}/\text{K}$ )	Measurement temperature ( $^{\circ}\text{C}$ )	Reference
0.3250	-10.8	37	[8]
0.4416	-11.6	37	[8]
0.4579	-12.0	20	[5]
0.6328	-12.4	20	[5]
	-12.0	37	[8, 9]
1.15	-12.6	20	[5]
	-12.7	37	[8, 9]
3.39	-12.4	20	[5]
	-13.0	37	[8, 9]
10.6	-9.8	20	[5]
$\infty$	265	7-47	[10]

**TABLE IV**  
**Room-Temperature Classical Oscillator Model Parameters**

Mode number $i$	$\nu_i$ ( $\text{cm}^{-1}$ )	$\Delta\epsilon_i$	$\Gamma_i/\nu_i$	Reference
1	217	4.0	0.0170	[16]
2	316	0.07	0.25	
$\epsilon_{\infty} = 2.07$				
Modified				
1	217	4.3	0.017	[see text]
2	360	0.6	0.25	
$\epsilon = 2.05$				