



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 8, 2026 – 08:32 AM UTC

PDB ID : 2BXE / pdb\_00002bx  
Title : Human serum albumin complexed with diflunisal  
Authors : Ghuman, J.; Zunszain, P.A.; Petitpas, I.; Bhattacharya, A.A.; Curry, S.  
Deposited on : 2005-07-26  
Resolution : 2.95 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at [validation@mail.wwpdb.org](mailto:validation@mail.wwpdb.org)

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4-5-2 with Phenix2.0  
Mogul : 2022.3.0, CSD as543be (2022)  
Xtriage (Phenix) : 2.0  
EDS : 3.0  
Buster-report : wwPDB partial adaption of 1.1.7 (2018)  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
CCP4 : 9.0.010 (Gargrove)  
Density-Fitness : 1.0.12  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

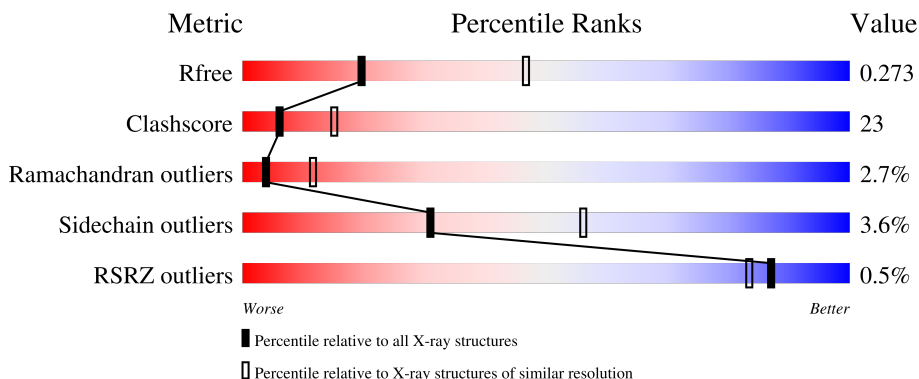
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.95 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	180053	1130 (2.98-2.94)
Clashscore	190562	1157 (2.98-2.94)
Ramachandran outliers	187476	1101 (2.98-2.94)
Sidechain outliers	187428	1101 (2.98-2.94)
RSRZ outliers	180081	1130 (2.98-2.94)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	585	
1	B	585	

## 2 Entry composition [i](#)

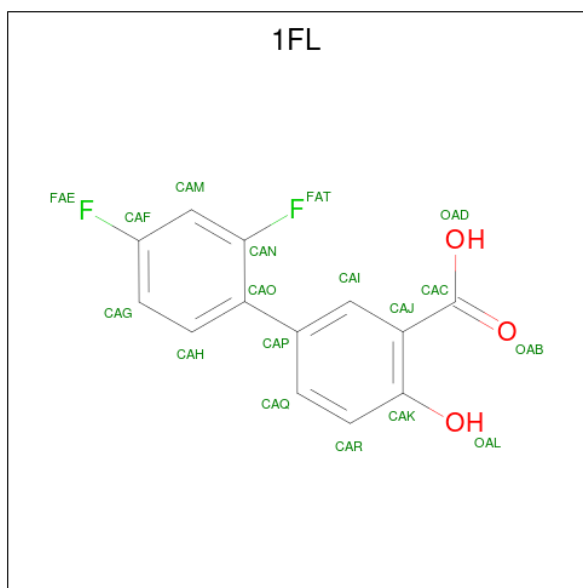
There are 2 unique types of molecules in this entry. The entry contains 8710 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

- Molecule 1 is a protein called SERUM ALBUMIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	569	Total	C	N	O	S	0	0	0
			4344	2743	722	839	40			
1	B	566	Total	C	N	O	S	0	0	0
			4258	2692	710	816	40			

- Molecule 2 is 5-(2,4-DIFLUOROPHENYL)-2-HYDROXY-BENZOIC ACID (CCD ID: 1FL) (formula: C<sub>13</sub>H<sub>8</sub>F<sub>2</sub>O<sub>3</sub>).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	F	O		
2	A	1	Total	C	F	O	0	0
			18	13	2	3		
2	A	1	Total	C	F	O	0	0
			18	13	2	3		
2	A	1	Total	C	F	O	0	0
			18	13	2	3		
2	B	1	Total	C	F	O	0	0
			18	13	2	3		

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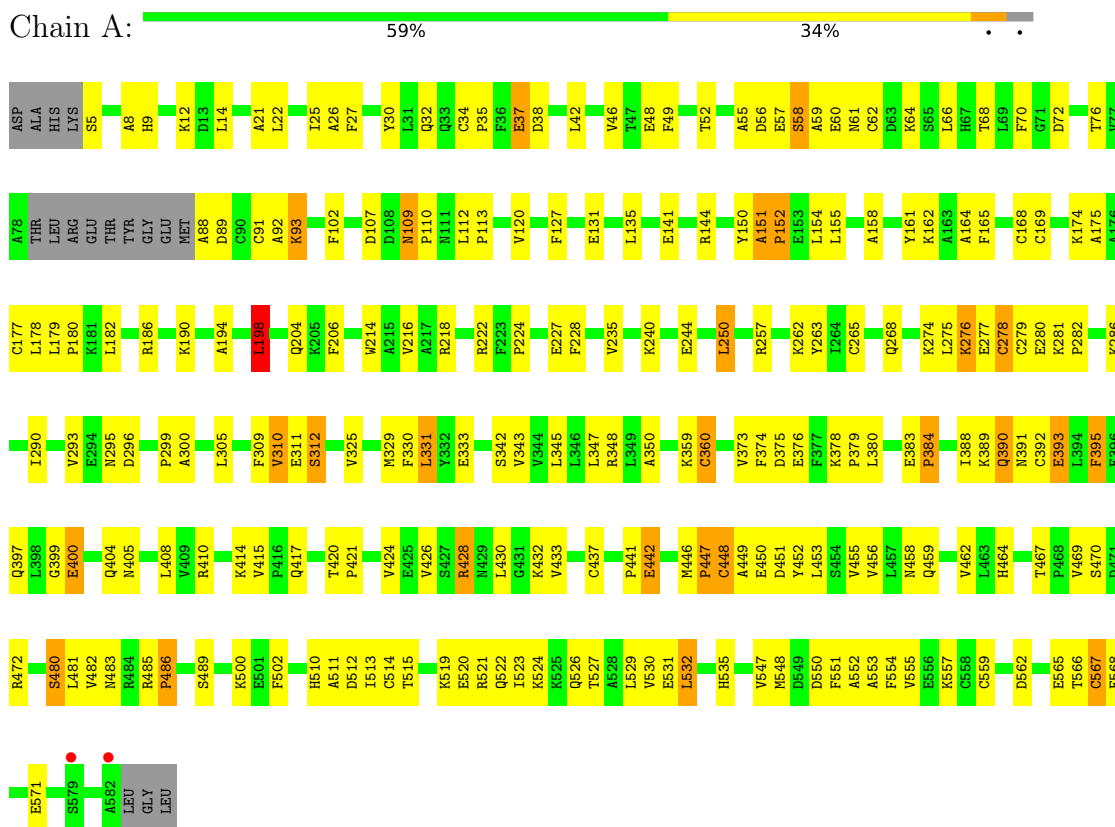
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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	F	O		
2	B	1	18	13	2	3	0	0
2	B	1	18	13	2	3	0	0

### 3 Residue-property plots [i](#)

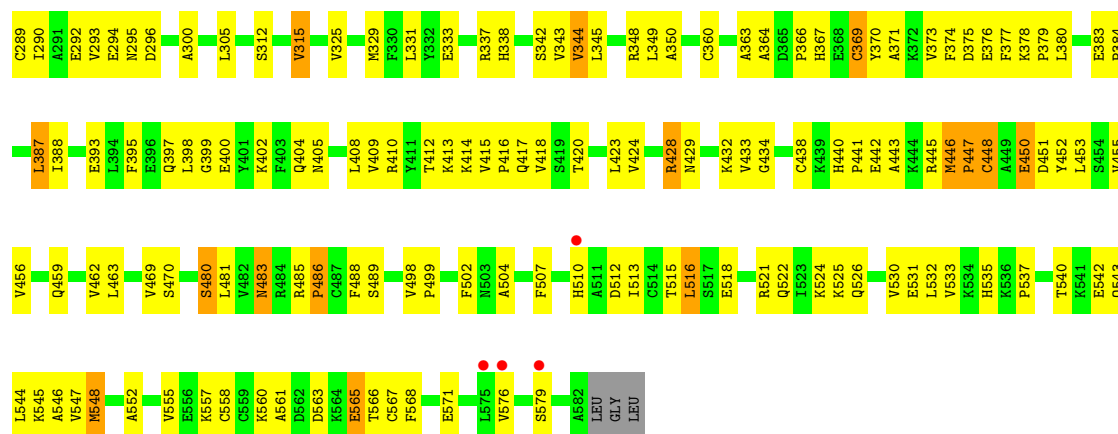
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

#### • Molecule 1: SERUM ALBUMIN



#### • Molecule 1: SERUM ALBUMIN





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	55.41Å 55.53Å 120.52Å 81.54° 90.09° 65.93°	Depositor
Resolution (Å)	38.22 – 2.95 38.22 – 2.95	Depositor EDS
% Data completeness (in resolution range)	98.4 (38.22-2.95) 98.4 (38.22-2.95)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.29 (at 2.95Å)	Xtrriage
Refinement program	CNS 1.1	Depositor
R, $R_{free}$	0.226 , 0.270 0.240 , 0.273	Depositor DCC
$R_{free}$ test set	1296 reflections (4.75%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	73.1	Xtrriage
Anisotropy	0.328	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.28 , 66.5	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.33$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	8710	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	82.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.66% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 1FL

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 5$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
1	A	0.56	0/4428	1.05	19/6011 (0.3%)
1	B	0.55	1/4339 (0.0%)	1.03	16/5895 (0.3%)
All	All	0.56	1/8767 (0.0%)	1.04	35/11906 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	325	VAL	CA-CB	-5.77	1.47	1.54

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	565	GLU	N-CA-C	-10.47	99.27	113.30
1	A	415	VAL	CA-C-N	8.09	128.85	119.47
1	A	415	VAL	C-N-CA	8.09	128.85	119.47
1	A	400	GLU	N-CA-C	7.58	119.18	111.07
1	B	338	HIS	N-CA-C	7.26	121.09	108.75
1	B	168	CYS	N-CA-C	7.15	120.11	111.82
1	B	450	GLU	N-CA-C	6.91	119.65	111.71
1	B	387	LEU	N-CA-C	-6.61	104.00	111.14
1	B	204	GLN	N-CA-C	6.57	118.10	111.07
1	A	449	ALA	N-CA-C	6.42	118.81	111.11
1	B	378	LYS	CA-C-N	6.33	125.56	118.97
1	B	378	LYS	C-N-CA	6.33	125.56	118.97
1	A	428	ARG	N-CA-C	-6.29	104.34	111.14
1	A	204	GLN	N-CA-C	6.11	118.02	111.36
1	A	395	PHE	N-CA-C	-6.08	104.28	111.03
1	B	250	LEU	N-CA-C	6.07	117.90	111.28
1	B	369	CYS	N-CA-C	6.00	117.48	111.07
1	A	250	LEU	N-CA-C	5.96	118.55	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	360	CYS	N-CA-C	5.92	117.81	111.36
1	B	428	ARG	N-CA-C	-5.89	104.78	111.14
1	B	415	VAL	CA-C-N	5.76	125.61	119.28
1	B	415	VAL	C-N-CA	5.76	125.61	119.28
1	A	567	CYS	N-CA-C	-5.72	105.41	112.38
1	B	393	GLU	N-CA-C	-5.52	105.27	111.28
1	B	399	GLY	N-CA-C	-5.41	106.81	111.95
1	A	437	CYS	N-CA-C	5.38	119.97	113.41
1	A	309	PHE	N-CA-C	5.32	120.78	113.97
1	A	198	LEU	N-CA-C	-5.28	105.42	111.07
1	A	278	CYS	N-CA-C	5.27	116.71	111.07
1	B	483	ASN	N-CA-C	5.26	119.70	113.28
1	A	375	ASP	N-CA-C	-5.13	106.17	112.90
1	B	32	GLN	N-CA-C	5.13	117.61	111.71
1	A	393	GLU	N-CA-C	-5.08	106.59	112.89
1	A	399	GLY	N-CA-C	-5.06	105.32	112.81
1	A	93	LYS	N-CA-C	5.01	116.90	109.69

There are no chirality outliers.

There are no planarity outliers.

## 5.2 Too-close contacts [\(i\)](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4344	0	4056	174	0
1	B	4258	0	3962	205	0
2	A	54	0	18	3	0
2	B	54	0	18	7	0
All	All	8710	0	8054	379	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 23.

All (379) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:VAL:HG22	1:A:452:TYR:HD2	1.06	1.10
1:A:433:VAL:HG22	1:A:452:TYR:CD2	1.91	1.05
1:B:433:VAL:HG22	1:B:452:TYR:HD2	1.30	0.95
1:B:433:VAL:HG22	1:B:452:TYR:CD2	2.02	0.93
1:B:222:ARG:HD3	2:B:2003:1FL:FAE	1.59	0.91
1:B:120:VAL:HG21	1:B:175:ALA:HA	1.53	0.88
1:A:222:ARG:HD3	2:A:2003:1FL:FAE	1.63	0.88
1:A:151:ALA:HB3	1:A:152:PRO:HD3	1.55	0.87
1:B:364:ALA:O	1:B:366:PRO:HD3	1.75	0.87
1:B:95:GLU:OE1	1:B:99:ASN:HB2	1.74	0.85
1:B:57:GLU:HG3	1:B:58:SER:H	1.42	0.85
1:B:567:CYS:SG	1:B:571:GLU:HB2	2.16	0.85
1:B:151:ALA:HB3	1:B:152:PRO:HD3	1.56	0.85
1:A:276:LYS:HD2	1:A:277:GLU:H	1.42	0.83
1:A:222:ARG:HG3	1:A:295:ASN:OD1	1.79	0.82
1:B:540:THR:HG23	1:B:543:GLN:H	1.42	0.81
1:A:120:VAL:HG11	1:A:174:LYS:HB2	1.61	0.81
1:B:222:ARG:HG3	1:B:295:ASN:OD1	1.79	0.81
1:A:417:GLN:OE1	1:A:417:GLN:N	2.13	0.80
1:A:424:VAL:O	1:A:428:ARG:HG3	1.82	0.80
1:B:383:GLU:HB3	1:B:384:PRO:HD3	1.60	0.80
1:A:135:LEU:HD11	1:A:162:LYS:HG3	1.64	0.79
1:A:390:GLN:O	1:A:393:GLU:N	2.13	0.79
1:B:281:LYS:HD2	1:B:285:GLU:HG2	1.65	0.79
1:A:383:GLU:HB3	1:A:384:PRO:HD3	1.66	0.78
1:A:550:ASP:O	1:A:554:PHE:HD1	1.66	0.78
1:A:182:LEU:O	1:A:186:ARG:HG2	1.84	0.78
1:A:547:VAL:HG12	1:A:551:PHE:HE1	1.46	0.77
1:A:61:ASN:HD22	1:A:64:LYS:HE3	1.50	0.77
1:B:345:LEU:HD22	1:B:446:MET:CE	2.15	0.77
1:B:446:MET:HB3	1:B:447:PRO:HD3	1.69	0.75
1:B:59:ALA:HB3	1:B:62:CYS:SG	2.27	0.74
1:A:107:ASP:OD2	1:A:110:PRO:HA	1.87	0.74
1:B:265:CYS:O	1:B:268:GLN:HG3	1.88	0.74
1:A:547:VAL:CG1	1:A:551:PHE:HE1	2.00	0.74
1:B:110:PRO:HB2	1:B:112:LEU:CD2	2.17	0.74
1:A:169:CYS:HA	1:A:174:LYS:HD3	1.71	0.73
1:A:414:LYS:O	1:A:472:ARG:NH1	2.23	0.72
1:A:345:LEU:HD23	1:A:446:MET:HE1	1.72	0.71
1:B:127:PHE:O	1:B:131:GLU:HB3	1.91	0.71
1:B:106:LYS:HD3	1:B:147:PRO:HB2	1.71	0.71
1:A:485:ARG:HB3	1:A:486:PRO:HD3	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:376:GLU:O	1:A:379:PRO:HD2	1.91	0.70
1:A:373:VAL:HG13	1:A:374:PHE:HD1	1.55	0.70
1:B:499:PRO:HB3	1:B:535:HIS:O	1.89	0.70
1:A:400:GLU:O	1:A:404:GLN:HG3	1.91	0.70
1:A:120:VAL:HG21	1:A:175:ALA:HA	1.75	0.69
1:B:194:ALA:HB1	1:B:455:VAL:CG1	2.22	0.69
1:B:214:TRP:CD1	1:B:343:VAL:HG11	2.27	0.69
1:B:285:GLU:O	1:B:289:CYS:HB2	1.93	0.69
1:B:561:ALA:HB3	1:B:563:ASP:OD1	1.93	0.69
1:B:483:ASN:O	1:B:486:PRO:HD2	1.92	0.68
1:B:206:PHE:CD2	1:B:481:LEU:HD22	2.29	0.68
1:A:276:LYS:CD	1:A:277:GLU:H	2.06	0.68
1:B:510:HIS:HA	1:B:568:PHE:CD2	2.28	0.68
1:A:151:ALA:HB3	1:A:152:PRO:CD	2.24	0.68
1:B:485:ARG:HB3	1:B:486:PRO:HD3	1.76	0.68
1:A:281:LYS:HB2	1:A:282:PRO:CD	2.25	0.67
1:B:32:GLN:NE2	1:B:110:PRO:HG2	2.10	0.67
1:B:260:LEU:O	1:B:260:LEU:HD23	1.94	0.67
1:A:35:PRO:HD2	1:A:38:ASP:OD2	1.95	0.67
1:A:305:LEU:HD21	1:A:333:GLU:HB3	1.77	0.66
1:B:110:PRO:HB2	1:B:112:LEU:HD22	1.76	0.66
1:B:10:ARG:O	1:B:14:LEU:HD23	1.95	0.66
1:B:344:VAL:HG22	1:B:451:ASP:OD1	1.96	0.65
1:B:441:PRO:C	1:B:443:ALA:H	2.04	0.65
1:A:66:LEU:O	1:A:70:PHE:HD1	1.78	0.65
1:A:529:LEU:HB2	1:A:548:MET:HE1	1.77	0.65
1:B:565:GLU:O	1:B:568:PHE:HB3	1.96	0.65
1:B:263:TYR:C	1:B:263:TYR:CD2	2.75	0.64
1:B:290:ILE:O	1:B:293:VAL:HG12	1.98	0.64
1:A:483:ASN:O	1:A:486:PRO:HD2	1.98	0.64
1:A:276:LYS:HD3	1:A:277:GLU:HG2	1.81	0.63
1:A:214:TRP:CD1	1:A:343:VAL:HG11	2.33	0.63
1:A:529:LEU:HD22	1:A:548:MET:CE	2.28	0.63
1:B:66:LEU:O	1:B:70:PHE:HD1	1.82	0.63
1:B:400:GLU:O	1:B:404:GLN:HG3	1.97	0.62
1:B:441:PRO:O	1:B:443:ALA:N	2.31	0.62
1:B:151:ALA:HB3	1:B:152:PRO:CD	2.28	0.62
1:B:417:GLN:H	1:B:417:GLN:CD	2.07	0.62
1:A:198:LEU:HD22	1:A:455:VAL:HA	1.82	0.61
1:B:558:CYS:SG	1:B:571:GLU:HB3	2.40	0.61
1:A:240:LYS:O	1:A:244:GLU:HG3	2.01	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:281:LYS:HB2	1:A:282:PRO:HD2	1.82	0.61
1:B:9:HIS:CD2	1:B:13:ASP:OD2	2.54	0.60
1:B:116:VAL:O	1:B:118:PRO:HD3	2.01	0.60
1:A:408:LEU:HD11	1:A:526:GLN:HB3	1.83	0.60
1:A:224:PRO:HB2	1:A:299:PRO:HD3	1.83	0.60
1:A:519:LYS:O	1:A:523:ILE:HG13	2.02	0.60
1:B:441:PRO:C	1:B:443:ALA:N	2.58	0.60
1:B:42:LEU:O	1:B:46:VAL:HG23	2.02	0.59
1:B:168:CYS:SG	1:B:177:CYS:C	2.84	0.59
1:A:150:TYR:OH	2:A:2003:1FL:OAD	2.19	0.59
1:A:426:VAL:O	1:A:430:LEU:HG	2.03	0.59
1:B:32:GLN:HE22	1:B:110:PRO:CG	2.16	0.59
1:B:179:LEU:HB2	1:B:180:PRO:HD3	1.85	0.59
1:A:373:VAL:HG13	1:A:374:PHE:CD1	2.38	0.58
1:B:29:GLN:HG2	1:B:143:ALA:O	2.03	0.58
1:B:57:GLU:O	1:B:58:SER:C	2.46	0.58
1:B:510:HIS:HA	1:B:568:PHE:HD2	1.66	0.58
1:A:222:ARG:C	1:A:224:PRO:HD3	2.29	0.58
1:A:378:LYS:HB3	1:A:379:PRO:HD3	1.86	0.58
1:B:120:VAL:HG21	1:B:175:ALA:CA	2.31	0.58
1:B:57:GLU:OE1	1:B:57:GLU:HA	2.03	0.58
1:A:450:GLU:O	1:A:450:GLU:HG3	2.03	0.57
1:B:513:ILE:HD12	1:B:555:VAL:HG23	1.85	0.57
1:B:424:VAL:O	1:B:428:ARG:HG3	2.04	0.57
1:B:518:GLU:O	1:B:522:GLN:HG3	2.04	0.57
1:B:420:THR:HG23	1:B:530:VAL:HG11	1.87	0.57
1:A:510:HIS:HA	1:A:568:PHE:CB	2.35	0.57
1:B:532:LEU:HD21	1:B:547:VAL:HG11	1.87	0.57
1:B:27:PHE:CD2	1:B:74:LEU:HD21	2.40	0.56
1:B:32:GLN:NE2	1:B:110:PRO:CG	2.67	0.56
1:B:117:ARG:NH2	1:B:182:LEU:HB3	2.19	0.56
1:B:424:VAL:HG23	1:B:530:VAL:HG21	1.86	0.56
1:A:194:ALA:HB1	1:A:455:VAL:CG1	2.36	0.56
1:B:348:ARG:NH2	1:B:450:GLU:OE2	2.30	0.56
1:B:405:ASN:O	1:B:409:VAL:HG23	2.04	0.56
1:A:5:SER:HA	1:A:62:CYS:O	2.05	0.56
1:B:502:PHE:HE2	1:B:507:PHE:HB2	1.70	0.56
1:A:9:HIS:C	1:A:9:HIS:HD1	2.14	0.56
1:B:150:TYR:OH	2:B:2003:1FL:OAB	2.22	0.56
1:B:376:GLU:O	1:B:379:PRO:HD2	2.06	0.56
1:B:432:LYS:O	1:B:433:VAL:C	2.47	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:513:ILE:CD1	1:B:555:VAL:HG23	2.36	0.55
1:B:543:GLN:O	1:B:546:ALA:HB3	2.06	0.55
1:B:107:ASP:HB3	1:B:110:PRO:HG3	1.87	0.55
1:B:260:LEU:HD22	2:B:2003:1FL:HAR	1.87	0.55
1:B:408:LEU:HD11	1:B:526:GLN:HB3	1.89	0.55
1:B:151:ALA:CB	1:B:152:PRO:HD3	2.35	0.55
1:A:345:LEU:HD23	1:A:446:MET:CE	2.35	0.55
1:B:446:MET:CB	1:B:447:PRO:HD3	2.37	0.55
1:A:325:VAL:O	1:A:329:MET:HG3	2.07	0.55
1:A:8:ALA:O	1:A:12:LYS:HG2	2.07	0.54
1:A:550:ASP:O	1:A:554:PHE:CD1	2.55	0.54
1:B:558:CYS:SG	1:B:571:GLU:CB	2.95	0.54
1:B:502:PHE:CE2	1:B:507:PHE:HB2	2.42	0.54
1:A:14:LEU:HD13	1:A:22:LEU:HD12	1.90	0.54
1:A:447:PRO:O	1:A:448:CYS:C	2.50	0.54
1:B:535:HIS:C	1:B:537:PRO:HD3	2.31	0.54
1:B:305:LEU:HD21	1:B:333:GLU:HB3	1.90	0.54
1:A:417:GLN:H	1:A:417:GLN:CD	2.08	0.54
1:A:380:LEU:O	1:A:384:PRO:HD2	2.08	0.53
1:A:547:VAL:CG1	1:A:551:PHE:CE1	2.88	0.53
1:A:551:PHE:O	1:A:552:ALA:C	2.50	0.53
1:B:117:ARG:HH22	1:B:182:LEU:HB3	1.73	0.53
1:A:89:ASP:O	1:A:92:ALA:HB3	2.09	0.53
1:A:276:LYS:CD	1:A:277:GLU:HG2	2.38	0.53
1:B:345:LEU:HD22	1:B:446:MET:HE2	1.89	0.53
1:A:331:LEU:HD13	1:A:350:ALA:HB2	1.90	0.53
1:B:331:LEU:HD13	1:B:350:ALA:HB2	1.91	0.53
1:B:417:GLN:CD	1:B:417:GLN:N	2.66	0.53
1:A:522:GLN:O	1:A:526:GLN:HG3	2.09	0.53
1:A:388:ILE:O	1:A:389:LYS:C	2.52	0.53
1:B:567:CYS:SG	1:B:571:GLU:CB	2.95	0.53
1:A:179:LEU:HB2	1:A:180:PRO:HD3	1.90	0.52
1:B:115:LEU:HG	1:B:145:ARG:CZ	2.39	0.52
1:B:516:LEU:HD12	1:B:516:LEU:H	1.74	0.52
1:B:32:GLN:HE22	1:B:110:PRO:HG3	1.74	0.52
1:B:440:HIS:HB3	1:B:441:PRO:CD	2.39	0.52
1:B:480:SER:OG	1:B:483:ASN:HB2	2.09	0.52
1:A:279:CYS:HA	1:A:286:LYS:CD	2.40	0.52
1:A:49:PHE:HD2	1:A:49:PHE:O	1.93	0.52
1:A:168:CYS:SG	1:A:177:CYS:C	2.93	0.52
1:A:464:HIS:HE1	1:A:470:SER:H	1.58	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:540:THR:HG22	1:B:543:GLN:HB2	1.91	0.52
1:A:151:ALA:CB	1:A:152:PRO:HD3	2.36	0.51
1:A:224:PRO:HD2	1:A:296:ASP:HB3	1.92	0.51
1:B:260:LEU:HD23	1:B:260:LEU:C	2.35	0.51
1:A:480:SER:OG	1:A:483:ASN:HB2	2.10	0.51
1:A:342:SER:HB2	1:A:451:ASP:OD2	2.10	0.51
1:B:25:ILE:O	1:B:29:GLN:HG3	2.10	0.51
1:A:88:ALA:O	1:A:89:ASP:C	2.53	0.51
1:A:161:TYR:O	1:A:164:ALA:HB3	2.11	0.51
1:B:26:ALA:HB2	1:B:250:LEU:HD12	1.92	0.51
1:A:383:GLU:HB3	1:A:384:PRO:CD	2.40	0.51
1:B:237:ASP:CB	1:B:260:LEU:HD12	2.41	0.51
1:B:51:LYS:HA	1:B:54:VAL:HG23	1.92	0.50
1:B:418:VAL:CG1	1:B:423:LEU:HG	2.41	0.50
1:B:408:LEU:O	1:B:412:THR:OG1	2.20	0.50
1:A:531:GLU:OE1	1:A:531:GLU:HA	2.12	0.50
1:A:49:PHE:C	1:A:49:PHE:CD2	2.88	0.50
1:A:464:HIS:CE1	1:A:469:VAL:H	2.30	0.50
1:A:343:VAL:O	1:A:347:LEU:HG	2.12	0.50
1:B:456:VAL:O	1:B:459:GLN:N	2.42	0.50
1:B:36:PHE:O	1:B:40:VAL:HG23	2.11	0.50
1:B:535:HIS:O	1:B:537:PRO:HD3	2.11	0.50
1:A:420:THR:O	1:A:424:VAL:HG23	2.11	0.50
1:B:409:VAL:O	1:B:413:LYS:HG3	2.12	0.49
1:B:150:TYR:HB2	1:B:196:GLN:HG2	1.94	0.49
1:B:224:PRO:HD2	1:B:296:ASP:HB3	1.95	0.49
1:A:186:ARG:O	1:A:190:LYS:HG3	2.12	0.49
1:B:384:PRO:O	1:B:388:ILE:HG12	2.13	0.49
1:B:373:VAL:HG13	1:B:374:PHE:HD1	1.76	0.49
1:A:42:LEU:O	1:A:46:VAL:HG23	2.12	0.49
1:B:100:GLU:HG2	1:B:103:LEU:HD12	1.93	0.49
1:A:49:PHE:O	1:A:49:PHE:CD2	2.66	0.49
1:B:222:ARG:HA	1:B:295:ASN:OD1	2.13	0.49
1:B:558:CYS:C	1:B:560:LYS:H	2.20	0.49
1:B:552:ALA:O	1:B:555:VAL:HG12	2.13	0.49
1:A:441:PRO:O	1:A:442:GLU:C	2.56	0.48
1:B:384:PRO:HB2	1:B:446:MET:HE1	1.94	0.48
1:A:512:ASP:O	1:A:515:THR:HG22	2.14	0.48
1:B:408:LEU:HD22	1:B:530:VAL:HG22	1.95	0.48
1:A:155:LEU:O	1:A:158:ALA:HB3	2.13	0.48
1:B:576:VAL:HA	1:B:579:SER:HB2	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:CYS:HA	1:A:286:LYS:HD2	1.96	0.48
1:A:342:SER:HB3	1:A:447:PRO:HA	1.94	0.48
1:B:66:LEU:O	1:B:70:PHE:CD1	2.64	0.48
1:A:165:PHE:CE1	1:A:178:LEU:HD21	2.48	0.48
1:B:481:LEU:N	2:B:2002:1FL:OAB	2.42	0.48
1:B:483:ASN:C	1:B:486:PRO:HD2	2.38	0.48
1:B:498:VAL:O	1:B:498:VAL:HG23	2.14	0.48
1:B:453:LEU:HD22	2:B:2001:1FL:HAQ	1.96	0.48
1:A:27:PHE:CE2	1:A:70:PHE:HD2	2.32	0.47
1:B:237:ASP:HB3	1:B:260:LEU:HD12	1.96	0.47
1:B:342:SER:HB3	1:B:447:PRO:HA	1.95	0.47
1:B:483:ASN:O	1:B:486:PRO:CD	2.62	0.47
1:A:26:ALA:HB2	1:A:250:LEU:HD12	1.96	0.47
1:A:275:LEU:HD23	1:A:293:VAL:HG11	1.96	0.47
1:B:274:LYS:CE	1:B:296:ASP:HA	2.44	0.47
1:A:66:LEU:O	1:A:70:PHE:CD1	2.65	0.47
1:A:141:GLU:OE1	1:A:144:ARG:HD3	2.14	0.47
1:A:345:LEU:CD2	1:A:446:MET:HE1	2.41	0.47
1:B:447:PRO:O	1:B:448:CYS:C	2.57	0.47
1:A:218:ARG:HD3	1:A:343:VAL:HG21	1.97	0.47
1:B:542:GLU:OE1	1:B:542:GLU:HA	2.14	0.47
1:A:550:ASP:O	1:A:553:ALA:HB3	2.15	0.47
1:B:453:LEU:CD2	2:B:2001:1FL:HAQ	2.44	0.47
1:B:9:HIS:HD2	1:B:13:ASP:OD2	1.96	0.47
1:B:540:THR:HG22	1:B:543:GLN:CG	2.45	0.47
1:A:32:GLN:HG2	1:A:144:ARG:O	2.15	0.47
1:A:432:LYS:O	1:A:433:VAL:C	2.56	0.47
1:A:532:LEU:O	1:A:535:HIS:HB3	2.15	0.47
1:B:521:ARG:O	1:B:525:LYS:HG3	2.15	0.46
1:A:348:ARG:HG3	1:A:482:VAL:HG12	1.96	0.46
1:B:384:PRO:HB2	1:B:446:MET:SD	2.55	0.46
1:B:151:ALA:CB	1:B:152:PRO:CD	2.93	0.46
1:B:398:LEU:HB3	1:B:402:LYS:HB2	1.96	0.46
1:B:433:VAL:O	1:B:434:GLY:C	2.54	0.46
1:B:558:CYS:C	1:B:560:LYS:N	2.71	0.46
1:B:138:TYR:O	1:B:142:ILE:HG12	2.14	0.46
1:B:51:LYS:C	1:B:53:CYS:H	2.23	0.46
1:A:390:GLN:O	1:A:392:CYS:N	2.49	0.46
1:A:566:THR:C	1:A:568:PHE:N	2.73	0.46
1:A:206:PHE:CZ	1:A:481:LEU:HB2	2.51	0.46
1:A:420:THR:HB	1:A:421:PRO:HD3	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:312:SER:O	1:B:315:VAL:HG23	2.16	0.46
1:A:127:PHE:CE1	1:A:131:GLU:HG3	2.51	0.45
1:A:265:CYS:O	1:A:268:GLN:HG3	2.15	0.45
1:B:223:PHE:CD1	1:B:272:SER:HB2	2.51	0.45
1:A:555:VAL:O	1:A:559:CYS:HB2	2.16	0.45
1:B:522:GLN:HA	1:B:525:LYS:HD3	1.98	0.45
1:B:462:VAL:HG23	1:B:463:LEU:N	2.32	0.45
1:A:206:PHE:CE2	1:A:481:LEU:HB2	2.50	0.45
1:B:292:GLU:O	1:B:292:GLU:HG2	2.17	0.45
1:B:395:PHE:C	1:B:397:GLN:N	2.74	0.45
1:A:61:ASN:ND2	1:A:64:LYS:HE3	2.27	0.45
1:B:521:ARG:O	1:B:524:LYS:HB2	2.16	0.45
1:A:420:THR:CG2	1:A:527:THR:HG23	2.47	0.45
1:A:456:VAL:O	1:A:459:GLN:N	2.50	0.45
1:B:370:TYR:CD1	1:B:370:TYR:C	2.95	0.45
1:A:21:ALA:HB1	1:A:155:LEU:HD21	1.99	0.45
1:A:311:GLU:O	1:A:312:SER:C	2.59	0.45
1:A:305:LEU:CD2	1:A:333:GLU:HB3	2.46	0.45
1:B:429:ASN:O	1:B:432:LYS:HB2	2.17	0.45
1:A:66:LEU:HD22	1:A:70:PHE:HE1	1.82	0.45
1:A:262:LYS:O	1:A:263:TYR:C	2.59	0.45
1:A:513:ILE:HD11	1:A:555:VAL:HG13	1.99	0.45
1:B:485:ARG:HB3	1:B:485:ARG:HE	1.63	0.45
1:A:520:GLU:O	1:A:524:LYS:HG3	2.17	0.45
1:B:19:PHE:O	1:B:23:VAL:HG23	2.17	0.45
1:A:483:ASN:C	1:A:486:PRO:HD2	2.42	0.44
1:A:500:LYS:O	1:A:535:HIS:ND1	2.47	0.44
1:A:532:LEU:HD11	1:A:547:VAL:HG11	1.99	0.44
1:A:290:ILE:HA	1:A:293:VAL:HG13	1.99	0.44
1:A:511:ALA:O	1:A:514:CYS:SG	2.76	0.44
1:A:405:ASN:O	1:A:408:LEU:HB2	2.17	0.44
1:B:274:LYS:HE2	1:B:296:ASP:HA	1.98	0.44
1:B:533:VAL:CG2	1:B:544:LEU:HD21	2.48	0.44
1:B:16:GLU:O	1:B:20:LYS:HG3	2.18	0.44
1:B:360:CYS:SG	1:B:369:CYS:C	3.00	0.44
1:A:521:ARG:O	1:A:524:LYS:N	2.48	0.44
1:A:30:TYR:CD1	1:A:102:PHE:HB3	2.53	0.44
1:B:274:LYS:HD2	1:B:294:GLU:OE2	2.18	0.44
1:A:56:ASP:C	1:A:58:SER:H	2.26	0.44
1:B:14:LEU:N	1:B:14:LEU:HD22	2.33	0.44
1:A:216:VAL:HG22	1:A:235:VAL:HG21	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:286:LYS:O	1:B:289:CYS:HB3	2.18	0.44
1:A:383:GLU:O	1:A:384:PRO:C	2.56	0.43
1:A:34:CYS:HA	1:A:35:PRO:HD3	1.88	0.43
1:A:278:CYS:O	1:A:281:LYS:HG2	2.18	0.43
1:B:239:THR:O	1:B:243:THR:OG1	2.28	0.43
1:A:91:CYS:C	1:A:93:LYS:H	2.26	0.43
1:A:567:CYS:O	1:A:571:GLU:N	2.46	0.43
1:B:383:GLU:HB3	1:B:384:PRO:CD	2.41	0.43
1:B:502:PHE:CZ	1:B:504:ALA:HA	2.54	0.43
1:A:141:GLU:OE1	1:A:141:GLU:HA	2.17	0.43
1:A:257:ARG:O	1:A:257:ARG:HG2	2.18	0.43
1:A:458:ASN:O	1:A:462:VAL:HG22	2.18	0.43
1:B:161:TYR:O	1:B:164:ALA:HB3	2.17	0.43
1:A:120:VAL:HG21	1:A:175:ALA:CA	2.45	0.43
1:B:19:PHE:CD2	1:B:19:PHE:C	2.97	0.43
1:B:533:VAL:HG22	1:B:544:LEU:HD21	2.01	0.43
1:B:567:CYS:O	1:B:571:GLU:N	2.50	0.43
1:A:68:THR:O	1:A:72:ASP:OD2	2.37	0.43
1:B:110:PRO:HB2	1:B:112:LEU:HD21	1.97	0.43
1:B:414:LYS:HE3	1:B:488:PHE:O	2.19	0.43
1:B:557:LYS:HB3	1:B:571:GLU:OE2	2.18	0.43
1:A:151:ALA:O	1:A:152:PRO:C	2.61	0.43
1:A:112:LEU:HA	1:A:113:PRO:HD3	1.91	0.43
1:A:206:PHE:CE2	1:A:481:LEU:HD13	2.54	0.43
1:A:329:MET:O	1:A:330:PHE:C	2.62	0.42
1:B:194:ALA:HB1	1:B:455:VAL:HG13	2.01	0.42
1:B:209:ARG:HH21	2:B:2002:1FL:HAH	1.84	0.42
1:B:410:ARG:O	1:B:414:LYS:HG3	2.20	0.42
1:B:333:GLU:O	1:B:337:ARG:HG3	2.19	0.42
1:A:263:TYR:C	1:A:263:TYR:CD2	2.97	0.42
1:A:395:PHE:C	1:A:397:GLN:N	2.77	0.42
1:B:113:PRO:O	1:B:114:ARG:C	2.63	0.42
1:B:272:SER:HB3	1:B:275:LEU:HG	2.01	0.42
1:B:531:GLU:OE1	1:B:531:GLU:HA	2.19	0.42
1:A:483:ASN:O	1:A:486:PRO:CD	2.64	0.42
1:B:275:LEU:O	1:B:276:LYS:C	2.63	0.42
1:B:395:PHE:C	1:B:397:GLN:H	2.26	0.42
1:B:89:ASP:O	1:B:92:ALA:HB3	2.18	0.42
1:B:215:ALA:O	1:B:216:VAL:C	2.63	0.42
1:B:434:GLY:HA2	1:B:438:CYS:SG	2.60	0.42
1:A:25:ILE:HD13	1:A:154:LEU:HD23	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:GLU:O	1:A:52:THR:HG23	2.20	0.42
1:A:227:GLU:O	1:A:228:PHE:C	2.61	0.41
1:A:464:HIS:CE1	1:A:470:SER:H	2.36	0.41
1:B:74:LEU:C	1:B:76:THR:H	2.28	0.41
1:B:107:ASP:O	1:B:110:PRO:HG3	2.20	0.41
1:B:367:HIS:O	1:B:371:ALA:HB2	2.19	0.41
1:B:563:ASP:OD2	1:B:567:CYS:N	2.50	0.41
1:A:109:ASN:HD22	1:A:109:ASN:HA	1.62	0.41
1:B:566:THR:O	1:B:567:CYS:C	2.64	0.41
1:A:178:LEU:O	1:A:179:LEU:C	2.63	0.41
1:A:384:PRO:HB2	1:A:446:MET:SD	2.60	0.41
1:B:282:PRO:O	1:B:283:LEU:C	2.64	0.41
1:B:293:VAL:HG22	1:B:294:GLU:N	2.34	0.41
1:B:512:ASP:O	1:B:515:THR:HG22	2.21	0.41
1:B:540:THR:CG2	1:B:543:GLN:HB2	2.50	0.41
1:B:117:ARG:O	1:B:117:ARG:HG3	2.21	0.41
1:B:545:LYS:HA	1:B:548:MET:HB2	2.01	0.41
1:A:120:VAL:CG1	1:A:174:LYS:HB2	2.41	0.41
1:A:290:ILE:O	1:A:293:VAL:HG22	2.21	0.41
1:A:373:VAL:O	1:A:376:GLU:HB2	2.21	0.41
1:B:540:THR:CG2	1:B:543:GLN:HG3	2.51	0.41
1:B:107:ASP:OD2	1:B:110:PRO:HA	2.20	0.41
1:B:349:LEU:HD22	1:B:377:PHE:CG	2.55	0.41
1:B:417:GLN:HB2	1:B:470:SER:HB2	2.03	0.41
1:A:395:PHE:O	1:A:397:GLN:N	2.54	0.41
1:B:24:LEU:HD13	1:B:43:VAL:HG21	2.01	0.41
1:B:91:CYS:C	1:B:93:LYS:N	2.76	0.41
1:A:158:ALA:O	1:A:161:TYR:HB3	2.21	0.41
1:A:274:LYS:HA	1:A:276:LYS:HE3	2.02	0.41
1:A:453:LEU:CD2	2:A:2001:1FL:HAQ	2.51	0.41
1:B:32:GLN:NE2	1:B:144:ARG:O	2.45	0.41
1:B:237:ASP:HB2	1:B:260:LEU:HD12	2.02	0.41
1:B:238:LEU:HD12	1:B:238:LEU:HA	1.87	0.41
1:B:412:THR:O	1:B:416:PRO:HG3	2.21	0.41
1:B:459:GLN:O	1:B:462:VAL:HG22	2.21	0.41
1:A:502:PHE:CD1	1:A:502:PHE:C	2.98	0.40
1:B:276:LYS:O	1:B:277:GLU:C	2.64	0.40
1:A:9:HIS:C	1:A:9:HIS:ND1	2.73	0.40
1:A:310:VAL:HG11	1:A:374:PHE:CE1	2.56	0.40
1:B:408:LEU:HD22	1:B:530:VAL:CG2	2.51	0.40
1:B:441:PRO:O	1:B:445:ARG:HG3	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:ASP:O	1:A:76:THR:HG23	2.20	0.40
1:A:174:LYS:O	1:A:175:ALA:C	2.63	0.40
1:A:359:LYS:HG3	1:A:360:CYS:N	2.36	0.40
1:A:410:ARG:O	1:A:414:LYS:HG3	2.22	0.40
1:B:61:ASN:HD22	1:B:64:LYS:CD	2.35	0.40
1:B:380:LEU:O	1:B:384:PRO:CD	2.68	0.40
1:A:37:GLU:CD	1:A:37:GLU:H	2.28	0.40
1:A:408:LEU:HD22	1:A:530:VAL:HG22	2.03	0.40
1:B:446:MET:CB	1:B:447:PRO:CD	2.99	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	565/585 (97%)	486 (86%)	65 (12%)	14 (2%)	4	13
1	B	562/585 (96%)	478 (85%)	68 (12%)	16 (3%)	4	11
All	All	1127/1170 (96%)	964 (86%)	133 (12%)	30 (3%)	4	11

All (30) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	59	ALA
1	A	60	GLU
1	B	54	VAL
1	B	57	GLU
1	B	58	SER
1	B	300	ALA
1	B	565	GLU
1	A	390	GLN
1	A	391	ASN

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Mol	Chain	Res	Type
1	A	562	ASP
1	B	442	GLU
1	A	58	SER
1	A	312	SER
1	A	447	PRO
1	A	557	LYS
1	B	315	VAL
1	B	363	ALA
1	A	300	ALA
1	A	448	CYS
1	B	447	PRO
1	A	57	GLU
1	B	276	LYS
1	A	55	ALA
1	B	60	GLU
1	B	75	CYS
1	B	281	LYS
1	B	448	CYS
1	B	118	PRO
1	B	469	VAL
1	A	151	ALA

### 5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	A	447/511 (88%)	432 (97%)	15 (3%)	32 58
1	B	432/511 (84%)	415 (96%)	17 (4%)	28 54
All	All	879/1022 (86%)	847 (96%)	32 (4%)	31 56

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	GLU
1	A	109	ASN

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<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	152	PRO
1	A	198	LEU
1	A	276	LYS
1	A	280	GLU
1	A	310	VAL
1	A	331	LEU
1	A	384	PRO
1	A	442	GLU
1	A	467	THR
1	A	480	SER
1	A	486	PRO
1	A	489	SER
1	A	532	LEU
1	B	112	LEU
1	B	152	PRO
1	B	167	GLU
1	B	182	LEU
1	B	222	ARG
1	B	245	CYS
1	B	263	TYR
1	B	329	MET
1	B	344	VAL
1	B	375	ASP
1	B	387	LEU
1	B	446	MET
1	B	480	SER
1	B	486	PRO
1	B	489	SER
1	B	516	LEU
1	B	548	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

<b>Mol</b>	<b>Chain</b>	<b>Res</b>	<b>Type</b>
1	A	29	GLN
1	A	39	HIS
1	A	61	ASN
1	A	99	ASN
1	A	109	ASN
1	A	268	GLN
1	A	338	HIS
1	A	429	ASN

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Mol	Chain	Res	Type
1	A	458	ASN
1	A	459	GLN
1	A	464	HIS
1	A	483	ASN
1	B	9	HIS
1	B	111	ASN
1	B	338	HIS
1	B	464	HIS
1	B	483	ASN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

## 5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

6 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with  $|Z| > 2$  is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	1FL	B	2002	-	19,19,19	1.16	1 (5%)	27,27,27	1.04	2 (7%)
2	1FL	A	2002	-	19,19,19	1.17	1 (5%)	27,27,27	1.08	1 (3%)
2	1FL	A	2001	-	19,19,19	1.23	2 (10%)	27,27,27	0.92	0
2	1FL	B	2003	-	19,19,19	1.30	2 (10%)	27,27,27	1.29	3 (11%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	1FL	A	2003	-	19,19,19	0.62	0	27,27,27	0.81	0
2	1FL	B	2001	-	19,19,19	1.22	1 (5%)	27,27,27	0.91	1 (3%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	1FL	B	2002	-	-	0/8/8/8	0/2/2/2
2	1FL	A	2002	-	-	0/8/8/8	0/2/2/2
2	1FL	A	2001	-	-	0/8/8/8	0/2/2/2
2	1FL	B	2003	-	-	0/8/8/8	0/2/2/2
2	1FL	A	2003	-	-	0/8/8/8	0/2/2/2
2	1FL	B	2001	-	-	0/8/8/8	0/2/2/2

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	2003	1FL	CAH-CAO	3.45	1.45	1.40
2	B	2001	1FL	CAO-CAN	2.74	1.43	1.39
2	A	2002	1FL	CAO-CAN	2.67	1.43	1.39
2	A	2001	1FL	CAO-CAN	2.64	1.43	1.39
2	B	2002	1FL	CAO-CAN	2.64	1.43	1.39
2	B	2003	1FL	CAH-CAG	2.52	1.42	1.38
2	A	2001	1FL	CAM-CAN	2.12	1.41	1.37

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	2003	1FL	CAN-CAM-CAF	2.79	119.64	116.67
2	A	2002	1FL	FAT-CAN-CAO	2.71	123.28	119.00
2	B	2002	1FL	FAT-CAN-CAO	2.68	123.24	119.00
2	B	2003	1FL	CAH-CAG-CAF	-2.43	115.88	118.38
2	B	2003	1FL	OAL-CAK-CAJ	2.15	125.70	121.71
2	B	2002	1FL	CAN-CAM-CAF	2.07	118.87	116.67
2	B	2001	1FL	FAT-CAN-CAO	2.05	122.23	119.00

There are no chirality outliers.

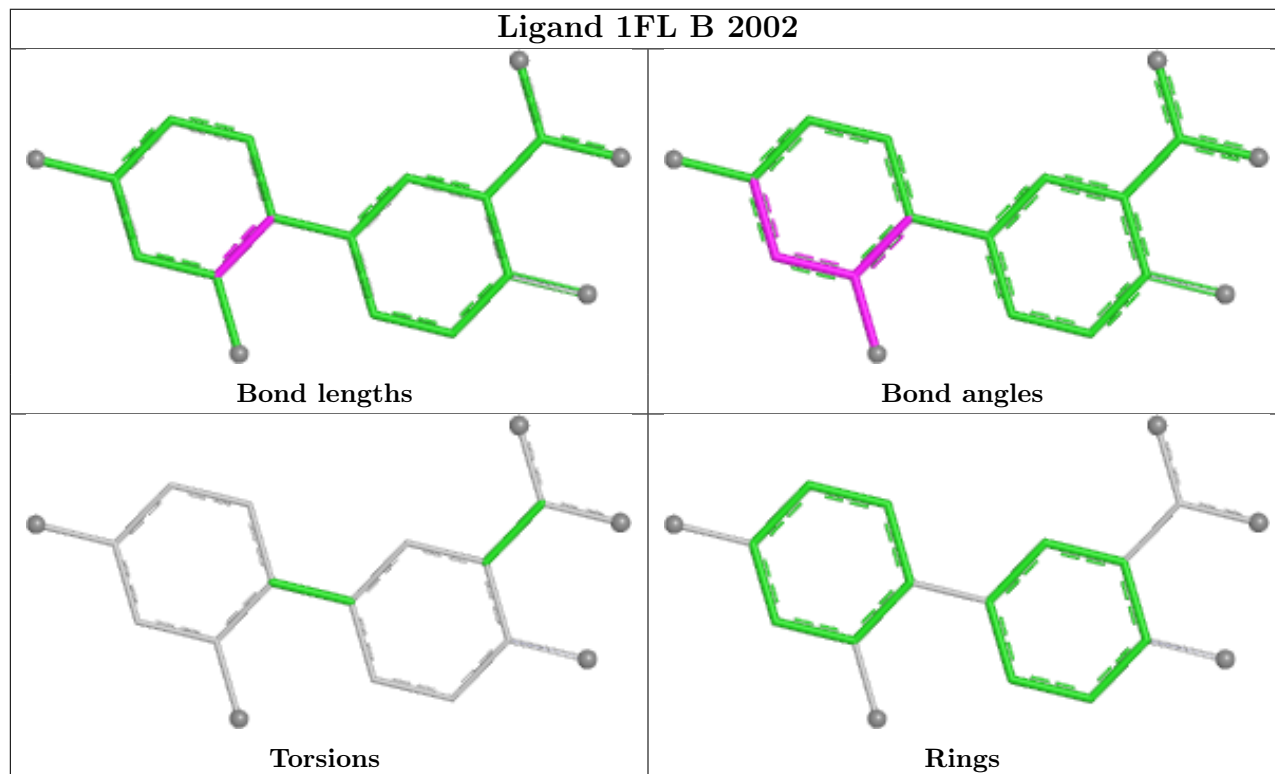
There are no torsion outliers.

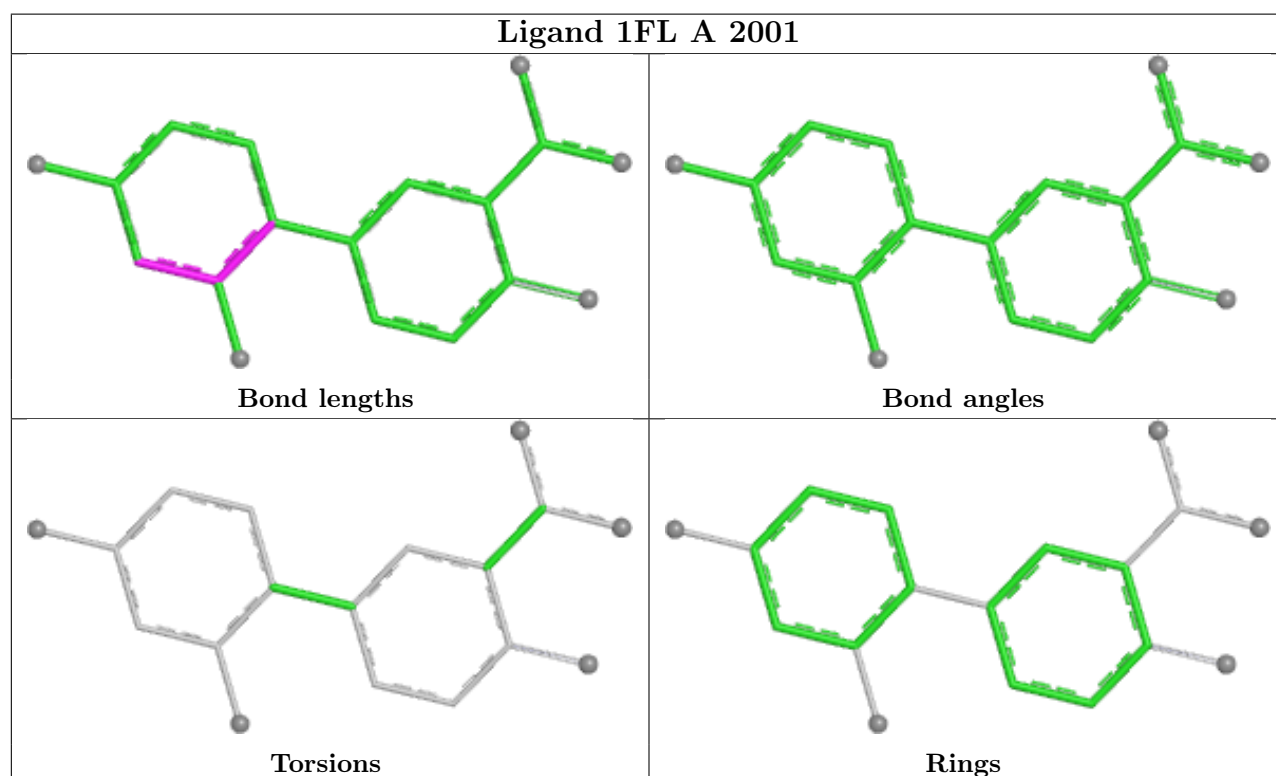
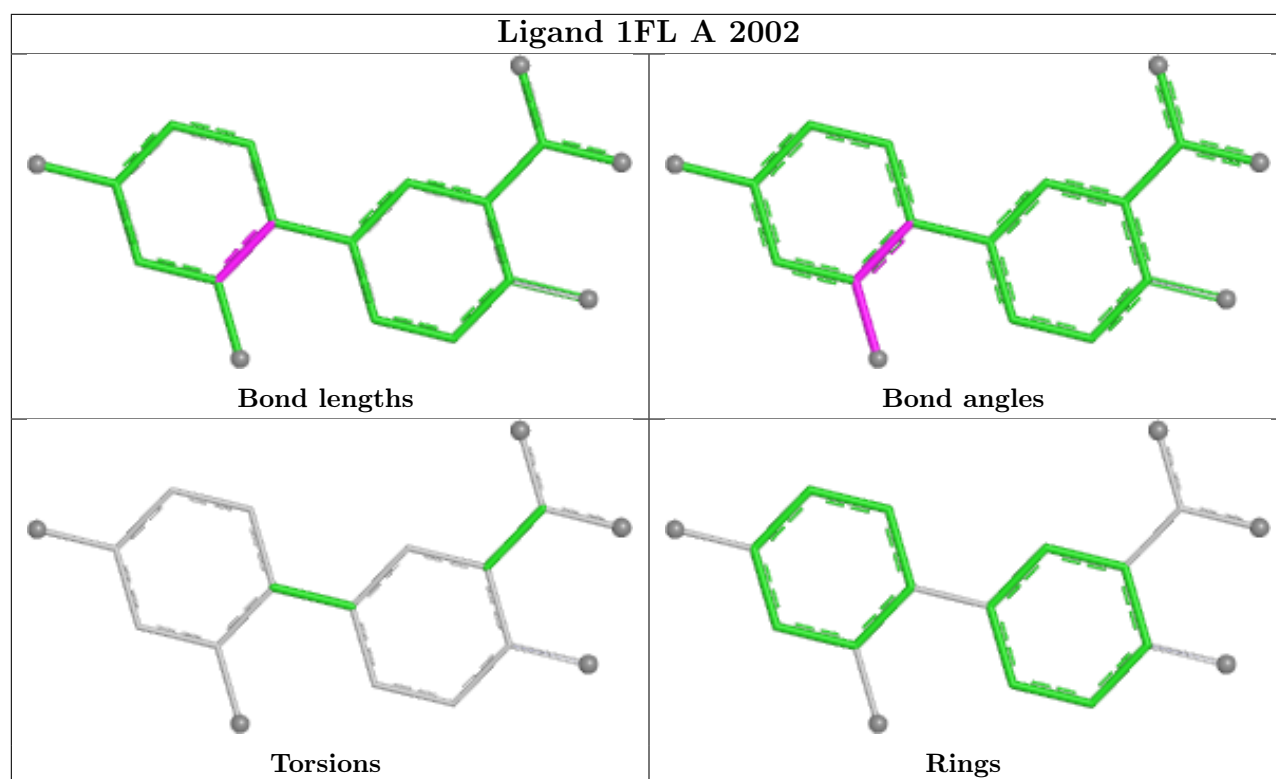
There are no ring outliers.

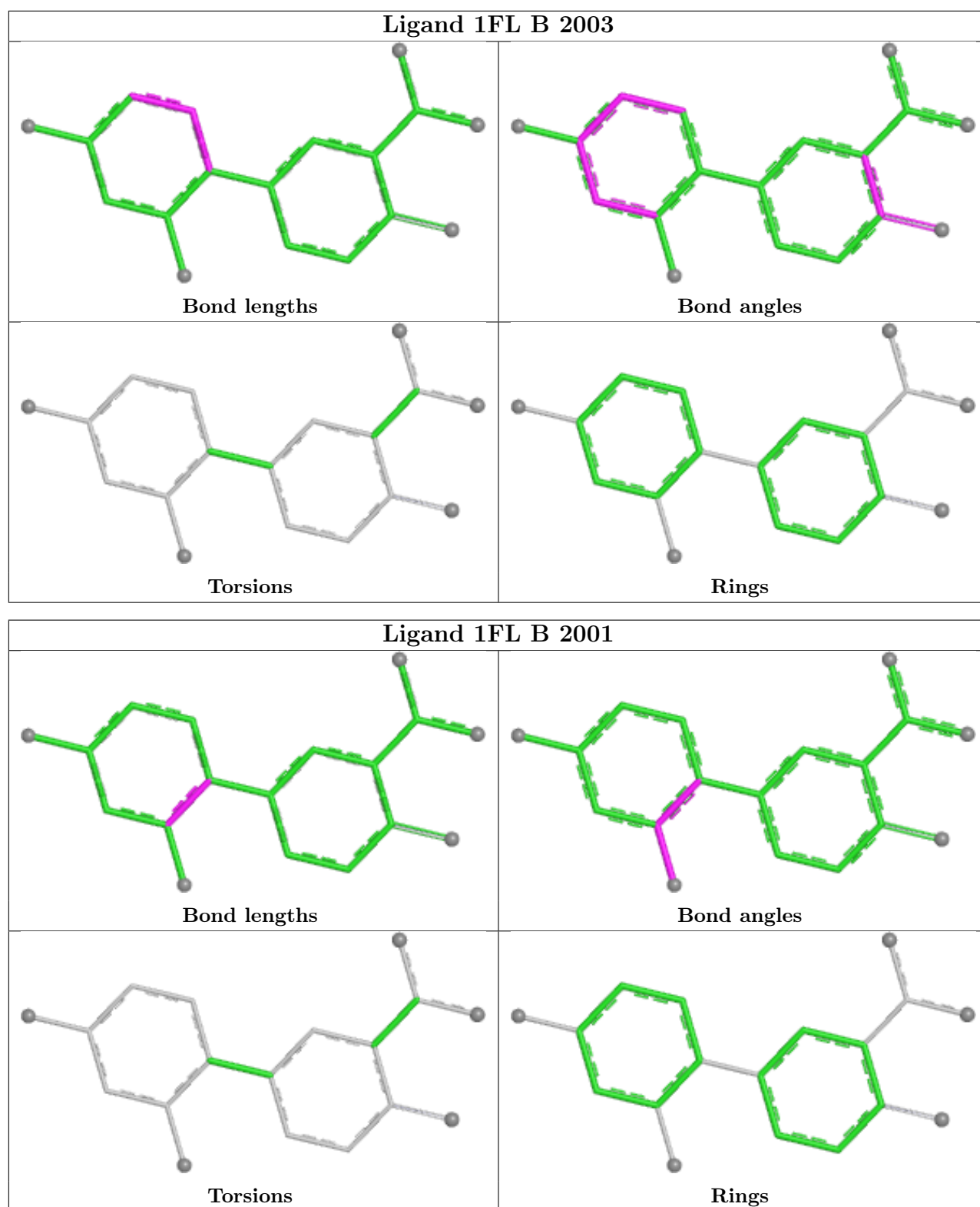
5 monomers are involved in 10 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	2002	1FL	2	0
2	A	2001	1FL	1	0
2	B	2003	1FL	3	0
2	A	2003	1FL	2	0
2	B	2001	1FL	2	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.







## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues

There are no chain breaks in this entry.

## 6 Fit of model and data [i](#)

### 6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å <sup>2</sup> )	Q<0.9
1	A	569/585 (97%)	-0.27	2 (0%) 88 86	35, 78, 156, 170	0
1	B	566/585 (96%)	-0.23	4 (0%) 84 80	35, 78, 148, 175	0
All	All	1135/1170 (97%)	-0.25	6 (0%) 87 83	35, 78, 150, 175	0

All (6) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	579	SER	4.2
1	A	579	SER	2.5
1	B	576	VAL	2.3
1	A	582	ALA	2.3
1	B	575	LEU	2.3
1	B	510	HIS	2.2

### 6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

### 6.3 Carbohydrates [i](#)

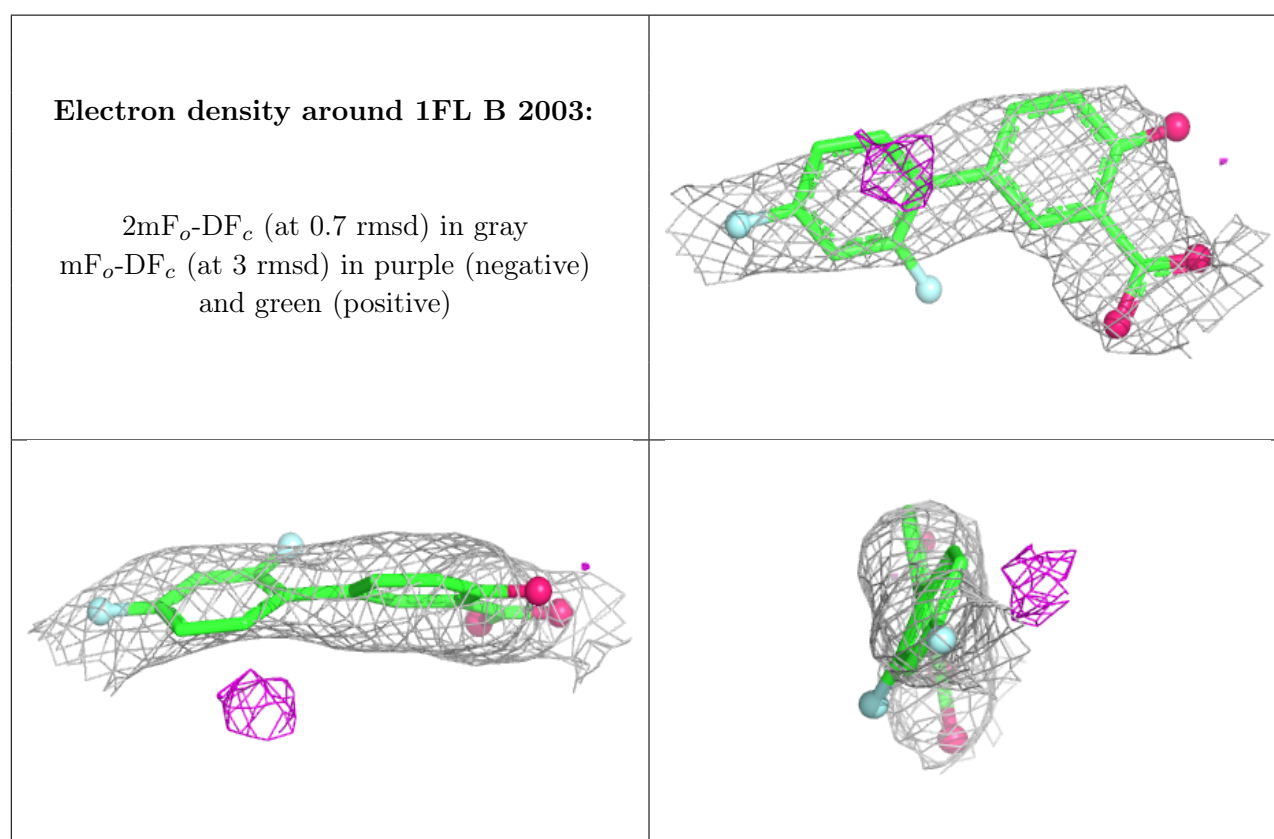
There are no oligosaccharides in this entry.

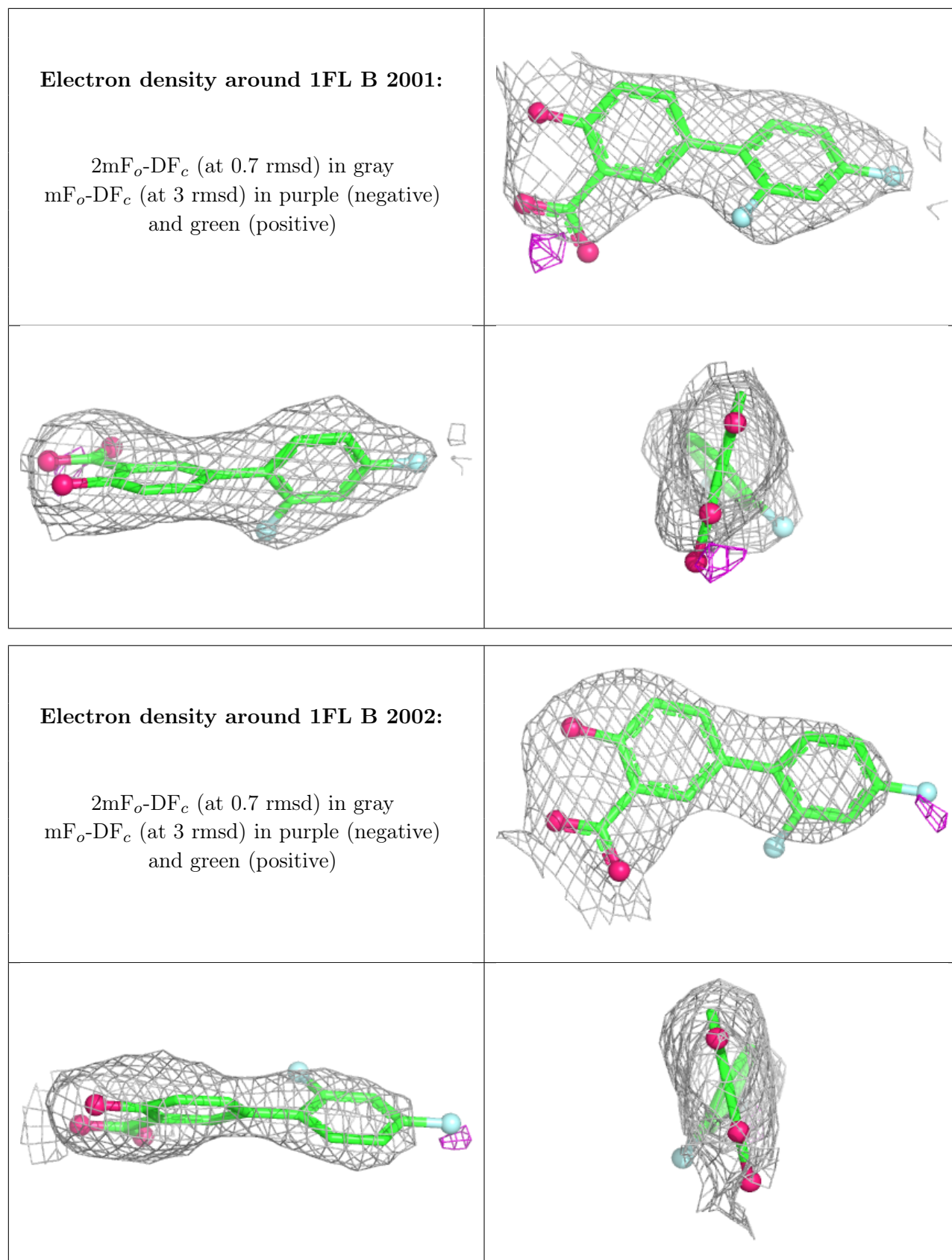
### 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors( $\text{\AA}^2$ )	Q<0.9
2	1FL	A	2003	18/18	0.90	0.13	109,109,111,111	0
2	1FL	B	2003	18/18	0.90	0.13	98,101,102,104	0
2	1FL	B	2001	18/18	0.91	0.13	63,67,71,73	0
2	1FL	B	2002	18/18	0.91	0.14	91,94,99,100	0
2	1FL	A	2002	18/18	0.91	0.12	77,82,91,92	0
2	1FL	A	2001	18/18	0.94	0.08	55,58,66,68	0

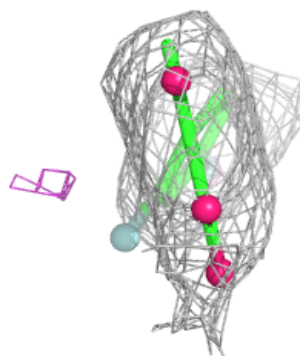
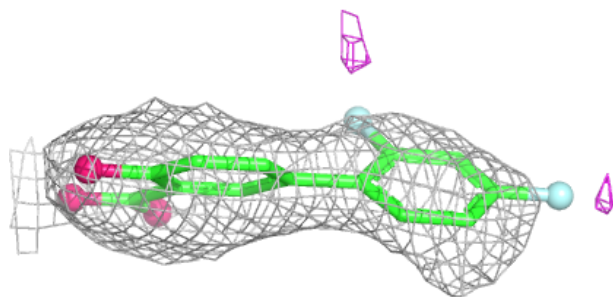
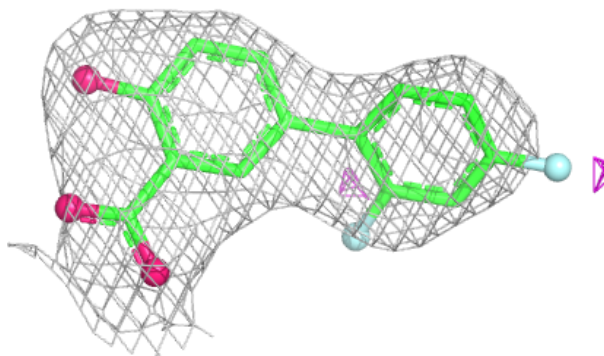
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



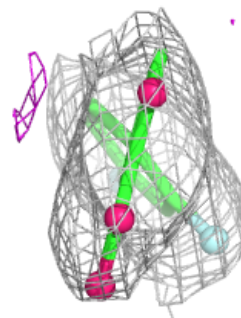
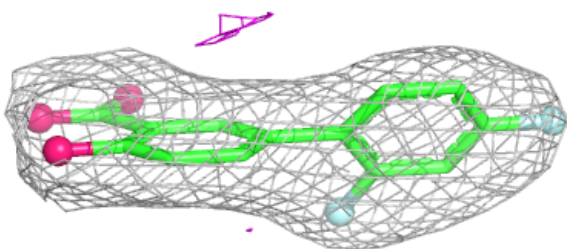
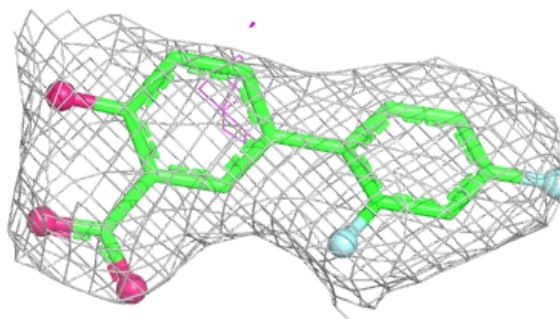


**Electron density around 1FL A 2002:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)

**Electron density around 1FL A 2001:**

$2mF_o-DF_c$  (at 0.7 rmsd) in gray  
 $mF_o-DF_c$  (at 3 rmsd) in purple (negative)  
and green (positive)



## 6.5 Other polymers [i](#)

There are no such residues in this entry.