



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 04:54 PM UTC

PDB ID : 3AVL / pdb_00003avl
Title : Crystal structures of novel allosteric peptide inhibitors of HIV integrase in the LEDGF binding site
Authors : Peat, T.S.; Deadman, J.J.; Newman, J.; Rhodes, D.I.
Deposited on : 2011-03-05
Resolution : 1.88 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

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>.

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
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

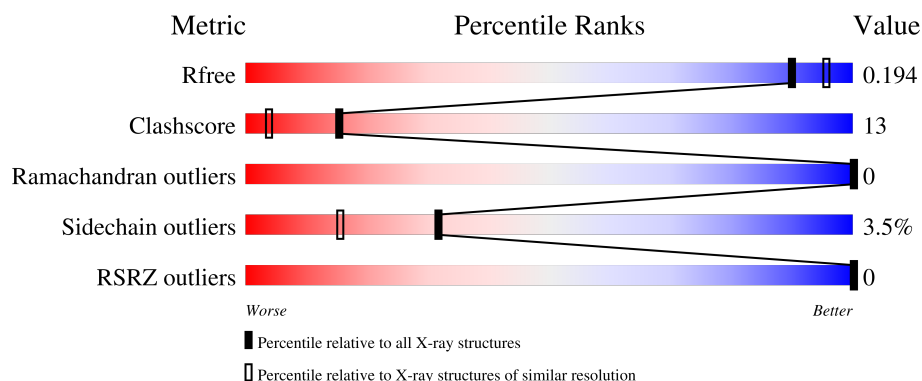
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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 ≥ 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	183	
1	B	183	
2	E	8	
2	F	8	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	A	4	-	-	X	-
3	SO4	A	7	-	-	X	-
3	SO4	B	3	-	-	X	-
3	SO4	B	8	-	-	X	-
5	ACY	A	1	-	-	X	-
5	ACY	A	214	-	-	X	-

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 2805 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 Integrase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	149	Total	C	N	O	S	0	11	0
			1230	788	210	228	4			
1	B	149	Total	C	N	O	S	0	13	0
			1238	794	211	229	4			

There are 46 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	MET	-	expression tag	UNP P12497
A	31	GLY	-	expression tag	UNP P12497
A	32	SER	-	expression tag	UNP P12497
A	33	SER	-	expression tag	UNP P12497
A	34	HIS	-	expression tag	UNP P12497
A	35	HIS	-	expression tag	UNP P12497
A	36	HIS	-	expression tag	UNP P12497
A	37	HIS	-	expression tag	UNP P12497
A	38	HIS	-	expression tag	UNP P12497
A	39	HIS	-	expression tag	UNP P12497
A	40	SER	-	expression tag	UNP P12497
A	41	SER	-	expression tag	UNP P12497
A	42	GLY	-	expression tag	UNP P12497
A	43	LEU	-	expression tag	UNP P12497
A	44	VAL	-	expression tag	UNP P12497
A	45	PRO	-	expression tag	UNP P12497
A	46	ARG	-	expression tag	UNP P12497
A	47	GLY	-	expression tag	UNP P12497
A	48	SER	-	expression tag	UNP P12497
A	49	HIS	-	expression tag	UNP P12497
A	56	SER	CYS	engineered mutation	UNP P12497
A	139	ASP	PHE	engineered mutation	UNP P12497
A	185	HIS	PHE	engineered mutation	UNP P12497
B	30	MET	-	expression tag	UNP P12497
B	31	GLY	-	expression tag	UNP P12497

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Chain	Residue	Modelled	Actual	Comment	Reference
B	32	SER	-	expression tag	UNP P12497
B	33	SER	-	expression tag	UNP P12497
B	34	HIS	-	expression tag	UNP P12497
B	35	HIS	-	expression tag	UNP P12497
B	36	HIS	-	expression tag	UNP P12497
B	37	HIS	-	expression tag	UNP P12497
B	38	HIS	-	expression tag	UNP P12497
B	39	HIS	-	expression tag	UNP P12497
B	40	SER	-	expression tag	UNP P12497
B	41	SER	-	expression tag	UNP P12497
B	42	GLY	-	expression tag	UNP P12497
B	43	LEU	-	expression tag	UNP P12497
B	44	VAL	-	expression tag	UNP P12497
B	45	PRO	-	expression tag	UNP P12497
B	46	ARG	-	expression tag	UNP P12497
B	47	GLY	-	expression tag	UNP P12497
B	48	SER	-	expression tag	UNP P12497
B	49	HIS	-	expression tag	UNP P12497
B	56	SER	CYS	engineered mutation	UNP P12497
B	139	ASP	PHE	engineered mutation	UNP P12497
B	185	HIS	PHE	engineered mutation	UNP P12497

- Molecule 2 is a protein called LEDGF peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	F	8	Total	C	N	O	0	1	0
			69	43	11	15			
2	E	8	Total	C	N	O	0	1	0
			69	43	11	15			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total	Cl	0	0
			1	1		
4	B	1	Total	Cl	0	0
			1	1		

- Molecule 5 is ACETIC ACID (CCD ID: ACY) (formula: C₂H₄O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			4	2	2		
5	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 6 is water.

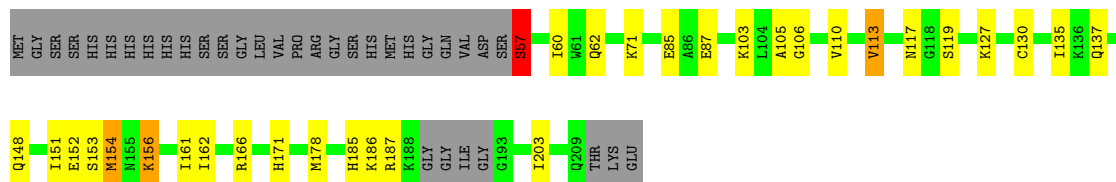
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	70	Total	O	0	0
			70	70		
6	B	76	Total	O	0	0
			76	76		
6	F	1	Total	O	0	0
			1	1		
6	E	2	Total	O	0	0
			2	2		

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.

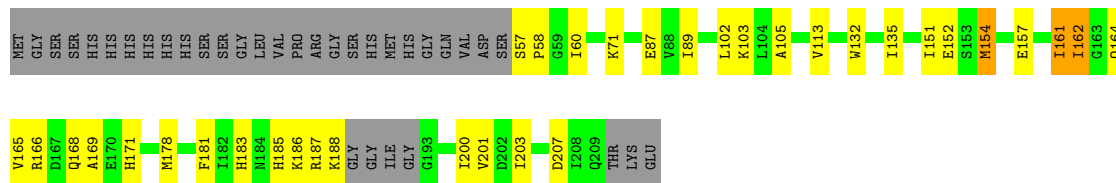
- Molecule 1: Integrase

Chain A: 



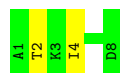
- Molecule 1: Integrase

Chain B: 



- Molecule 2: LEDGF peptide

Chain F: 



- Molecule 2: LEDGF peptide

Chain E: 



4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	70.48Å 70.48Å 66.86Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	61.00 – 1.88 61.00 – 1.88	Depositor EDS
% Data completeness (in resolution range)	100.0 (61.00-1.88) 100.0 (61.00-1.88)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.56 (at 1.88Å)	Xtriage
Refinement program	REFMAC 5.6.0062	Depositor
R, R_{free}	0.153 , 0.194 0.152 , 0.194	Depositor DCC
R_{free} test set	1531 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å ²)	21.1	Xtriage
Anisotropy	0.070	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 34.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.026 for -h,-k,l 0.487 for h,-h-k,-l 0.025 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	2805	wwPDB-VP
Average B, all atoms (Å ²)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.20% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical 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: ACY, CL, SO4

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	1.51	6/1275 (0.5%)	1.28	6/1726 (0.3%)
1	B	1.52	9/1286 (0.7%)	1.27	2/1742 (0.1%)
2	E	1.12	0/68	0.88	0/91
2	F	1.33	0/68	0.95	0/91
All	All	1.50	15/2697 (0.6%)	1.26	8/3650 (0.2%)

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	161	ILE	CA-CB	9.24	1.65	1.54
1	B	89	ILE	N-CA	6.57	1.52	1.46
1	A	203	ILE	CA-CB	6.45	1.61	1.54
1	A	161	ILE	CA-CB	6.17	1.62	1.54
1	B	169	ALA	N-CA	5.91	1.53	1.46
1	B	200	ILE	N-CA	5.62	1.53	1.46
1	B	203	ILE	CA-CB	5.46	1.60	1.54
1	B	132	TRP	C-O	5.45	1.30	1.24
1	A	110	VAL	CA-C	5.33	1.59	1.53
1	A	113[A]	VAL	C-O	5.28	1.29	1.24
1	A	113[B]	VAL	C-O	5.28	1.29	1.24
1	B	162	ILE	N-CA	-5.18	1.40	1.46
1	B	201	VAL	CA-CB	5.15	1.60	1.54
1	B	165	VAL	CA-CB	5.13	1.60	1.54
1	A	117	ASN	N-CA	5.01	1.51	1.46

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	57	SER	CA-C-N	-7.77	111.64	119.56
1	A	57	SER	C-N-CA	-7.77	111.64	119.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	130	CYS	CA-CB-SG	-5.61	101.50	114.40
1	A	119	SER	N-CA-C	-5.54	105.32	111.36
1	B	102	LEU	N-CA-C	-5.25	105.64	111.36
1	A	154	MET	CB-CG-SD	-5.21	97.05	112.70
1	A	85	GLU	N-CA-C	-5.21	101.27	109.50
1	B	154	MET	CB-CG-SD	-5.03	97.61	112.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

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	1230	0	1254	35	0
1	B	1238	0	1268	37	0
2	E	69	0	71	7	0
2	F	69	0	71	5	0
3	A	20	0	0	5	0
3	B	20	0	0	6	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
5	A	8	0	6	8	0
6	A	70	0	0	6	0
6	B	76	0	0	7	1
6	E	2	0	0	0	0
6	F	1	0	0	0	0
All	All	2805	0	2670	72	1

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 13.

All (72) 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)
3:B:6:SO4:O2	6:B:270:HOH:O	1.58	1.19

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:A:5:SO4:O1	6:A:266:HOH:O	1.65	1.11
1:A:185:HIS:CE1	5:A:1:ACY:H2	1.89	1.07
1:A:87[B]:GLU:OE1	6:A:239:HOH:O	1.81	0.97
1:B:87[B]:GLU:OE1	6:B:247:HOH:O	1.83	0.94
1:A:87[B]:GLU:OE2	1:B:103[B]:LYS:HE3	1.76	0.85
1:A:185:HIS:CE1	5:A:1:ACY:CH3	2.59	0.85
1:A:62[B]:GLN:NE2	1:A:151:ILE:HD11	1.93	0.83
1:A:62[B]:GLN:HE21	1:A:151:ILE:HD11	1.45	0.82
1:A:178:MET:CE	2:F:4[B]:ILE:HG21	2.11	0.81
1:B:178:MET:CE	2:E:4[B]:ILE:HG21	2.15	0.76
5:A:214:ACY:H3	1:B:185:HIS:NE2	2.02	0.74
3:B:1:SO4:O3	6:B:269:HOH:O	2.07	0.72
1:B:178:MET:HE1	2:E:4[B]:ILE:HG21	1.73	0.71
1:B:57[B]:SER:HA	1:B:60[B]:ILE:HG12	1.73	0.71
1:A:57:SER:HA	1:A:60[B]:ILE:HG12	1.73	0.70
1:A:103[B]:LYS:HE3	1:B:87[B]:GLU:OE2	1.92	0.69
1:A:178:MET:HE3	2:F:4[B]:ILE:HG21	1.76	0.68
1:B:71:LYS:NZ	6:B:237:HOH:O	2.27	0.67
3:A:7:SO4:O1	6:A:265:HOH:O	2.10	0.67
3:B:8:SO4:O2	6:B:268:HOH:O	2.13	0.65
1:A:178:MET:HE1	2:F:4[B]:ILE:HG21	1.80	0.64
1:A:71:LYS:NZ	6:A:236:HOH:O	2.31	0.64
3:A:4:SO4:O4	6:A:220:HOH:O	2.15	0.62
1:A:178:MET:HE3	2:F:4[A]:ILE:HG21	1.81	0.62
1:B:181:PHE:O	1:B:185:HIS:HD2	1.83	0.61
1:B:171:HIS:HD2	3:B:3:SO4:O1	1.86	0.59
1:A:103[B]:LYS:HZ3	1:B:103[B]:LYS:HZ1	1.50	0.59
1:A:171:HIS:HD2	3:A:4:SO4:O3	1.86	0.59
1:A:178:MET:CE	2:F:4[A]:ILE:HG21	2.33	0.59
1:A:103[B]:LYS:CE	1:B:87[B]:GLU:OE2	2.53	0.56
1:A:103[B]:LYS:NZ	1:B:103[B]:LYS:HZ1	2.05	0.55
1:A:62[B]:GLN:NE2	6:A:249:HOH:O	2.35	0.54
1:A:87[B]:GLU:OE2	1:B:103[B]:LYS:CE	2.53	0.53
1:B:57[B]:SER:HA	1:B:60[B]:ILE:CG1	2.37	0.53
1:A:57:SER:HA	1:A:60[B]:ILE:CG1	2.37	0.52
1:B:57[A]:SER:HB3	1:B:58:PRO:HD3	1.92	0.51
1:A:103[B]:LYS:NZ	1:B:103[B]:LYS:NZ	2.59	0.51
5:A:214:ACY:H3	1:B:185:HIS:CE1	2.46	0.50
1:B:87[B]:GLU:OE2	3:B:8:SO4:O3	2.29	0.50
1:A:87[B]:GLU:OE2	3:A:7:SO4:O3	2.31	0.48
5:A:214:ACY:H3	1:B:185:HIS:HE2	1.76	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:181:PHE:O	1:B:185:HIS:CD2	2.66	0.48
1:B:178:MET:HE3	2:E:4[B]:ILE:HG21	1.93	0.48
1:A:162:ILE:O	1:A:166:ARG:HB2	2.14	0.47
1:B:178:MET:HE3	2:E:4[A]:ILE:HG21	1.98	0.46
1:A:187:ARG:HH11	1:A:187:ARG:HG2	1.81	0.45
1:B:188:LYS:NZ	6:B:275:HOH:O	2.42	0.45
1:A:185:HIS:CE1	5:A:1:ACY:H1	2.49	0.45
1:A:127:LYS:NZ	1:A:137:GLN:HE22	2.15	0.44
1:B:162:ILE:O	1:B:166:ARG:HB2	2.17	0.44
1:B:171:HIS:CD2	3:B:3:SO4:O1	2.70	0.44
1:A:148:GLN:O	1:A:152[A]:GLU:HG3	2.18	0.44
1:A:187:ARG:HG2	1:A:187:ARG:NH1	2.33	0.43
1:B:161:ILE:HA	1:B:164:GLN:HE21	1.83	0.43
1:A:105:ALA:HB2	1:A:135:ILE:HD11	2.00	0.43
1:A:113[A]:VAL:HG12	1:A:135:ILE:HG21	2.01	0.43
1:B:187:ARG:HG2	1:B:187:ARG:HH11	1.83	0.43
1:B:168:GLN:O	2:E:4[B]:ILE:HG22	2.20	0.42
1:B:178:MET:CE	2:E:4[A]:ILE:HG21	2.49	0.42
1:B:187:ARG:HG2	1:B:187:ARG:NH1	2.35	0.42
1:A:106:GLY:O	5:A:214:ACY:H1	2.20	0.42
1:B:183:HIS:CE1	6:B:275:HOH:O	2.72	0.42
1:B:105:ALA:HB2	1:B:135:ILE:HD11	2.00	0.41
1:B:186:LYS:HD3	1:B:188:LYS:HE3	2.03	0.41
1:A:156:LYS:HB3	1:A:156:LYS:HE2	1.72	0.41
1:B:58:PRO:HG3	1:B:207:ASP:CG	2.46	0.41
1:B:154:MET:HB3	1:B:154:MET:HE2	1.81	0.41
2:E:6:ASN:OD1	2:E:8:ASP:HB2	2.20	0.41
1:A:106:GLY:C	5:A:214:ACY:H1	2.46	0.40
1:A:154:MET:HE2	1:A:154:MET:HB3	2.00	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
6:B:217:HOH:O	6:B:269:HOH:O[2_555]	1.99	0.21

5.3 Torsion angles

5.3.1 Protein backbone

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	156/183 (85%)	155 (99%)	1 (1%)	0	100	100
1	B	157/183 (86%)	157 (100%)	0	0	100	100
2	E	7/8 (88%)	6 (86%)	1 (14%)	0	100	100
2	F	7/8 (88%)	6 (86%)	1 (14%)	0	100	100
All	All	327/382 (86%)	324 (99%)	3 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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	134/150 (89%)	130 (97%)	4 (3%)	36	19
1	B	136/150 (91%)	132 (97%)	4 (3%)	37	20
2	E	8/7 (114%)	7 (88%)	1 (12%)	4	0
2	F	8/7 (114%)	7 (88%)	1 (12%)	4	0
All	All	286/314 (91%)	276 (96%)	10 (4%)	32	15

All (10) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	57	SER
1	A	153	SER

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Mol	Chain	Res	Type
1	A	156	LYS
1	A	186	LYS
1	B	151[A]	ILE
1	B	151[B]	ILE
1	B	152	GLU
1	B	157	GLU
2	F	2	THR
2	E	2	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	137	GLN
1	A	171	HIS
1	A	209	GLN
1	B	146	GLN
1	B	164	GLN
1	B	171	HIS
1	B	185	HIS
1	B	209	GLN

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](#)

Of 12 ligands modelled in this entry, 2 are monoatomic - leaving 10 for Mogul analysis.

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	$\# Z > 2$
3	SO4	A	2	-	4,4,4	0.12	0	6,6,6	1.48	1 (16%)
3	SO4	A	7	-	4,4,4	0.36	0	6,6,6	0.69	0
5	ACY	A	1	-	3,3,3	0.98	0	3,3,3	1.26	0
3	SO4	B	3	-	4,4,4	0.46	0	6,6,6	0.42	0
3	SO4	B	8	-	4,4,4	0.27	0	6,6,6	0.63	0
3	SO4	B	6	-	4,4,4	0.33	0	6,6,6	2.00	1 (16%)
3	SO4	A	4	-	4,4,4	0.20	0	6,6,6	0.70	0
3	SO4	B	1	-	4,4,4	0.46	0	6,6,6	0.88	0
3	SO4	A	5	-	4,4,4	0.44	0	6,6,6	1.84	2 (33%)
5	ACY	A	214	-	3,3,3	0.65	0	3,3,3	1.47	0

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	6	SO4	O4-S-O2	-3.53	91.11	109.56
3	A	5	SO4	O4-S-O1	3.16	126.07	109.56
3	A	5	SO4	O2-S-O1	-2.98	88.06	109.06
3	A	2	SO4	O4-S-O2	-2.76	95.14	109.56

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

9 monomers are involved in 19 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	7	SO4	2	0
5	A	1	ACY	3	0
3	B	3	SO4	2	0
3	B	8	SO4	2	0
3	B	6	SO4	1	0
3	A	4	SO4	2	0
3	B	1	SO4	1	0
3	A	5	SO4	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	214	ACY	5	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

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, 95th 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 ²)	Q<0.9
1	A	149/183 (81%)	-1.21	0 100 100	8, 18, 37, 46	28 (18%)
1	B	149/183 (81%)	-1.23	0 100 100	8, 18, 35, 51	30 (20%)
2	E	8/8 (100%)	-0.66	0 100 100	20, 41, 62, 71	1 (12%)
2	F	8/8 (100%)	-0.72	0 100 100	20, 40, 62, 69	1 (12%)
All	All	314/382 (82%)	-1.19	0 100 100	8, 18, 41, 71	60 (19%)

There are no RSRZ outliers to report.

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, 95th 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(Å ²)	Q<0.9
3	SO4	A	7	5/5	0.98	0.05	34,41,51,57	5
3	SO4	B	8	5/5	0.98	0.07	33,40,48,59	5
5	ACY	A	1	4/4	0.98	0.07	43,47,54,58	0
5	ACY	A	214	4/4	0.98	0.06	44,46,47,53	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	SO4	B	1	5/5	0.99	0.03	31,40,45,46	5
3	SO4	B	3	5/5	0.99	0.04	32,38,44,49	5
3	SO4	B	6	5/5	0.99	0.04	27,28,41,44	5
3	SO4	A	4	5/5	0.99	0.04	28,30,38,38	5
4	CL	A	213	1/1	0.99	0.05	48,48,48,48	0
4	CL	B	213	1/1	0.99	0.05	48,48,48,48	0
3	SO4	A	5	5/5	0.99	0.03	24,24,38,41	5
3	SO4	A	2	5/5	0.99	0.04	33,39,42,47	5

6.5 Other polymers [i](#)

There are no such residues in this entry.