



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 12:27 PM GMT

PDB ID : 4IOF
Title : Crystal structure analysis of Fab-bound human Insulin Degrading Enzyme (IDE)
Authors : McCord, L.A.; Liang, W.G; Hoey, R.; Dowdell, E.; Koide, A.; Koide, S.; Tang, W.J.
Deposited on : 2013-01-07
Resolution : 3.35 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

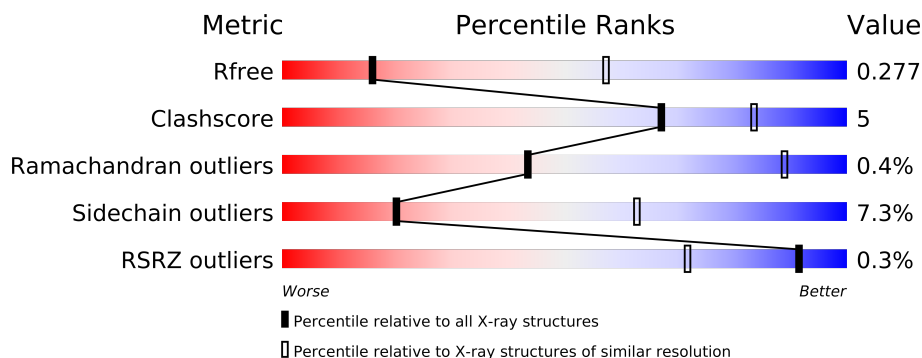
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.35 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}	66092	1141 (3.50-3.22)
Clashscore	79885	1030 (3.48-3.24)
Ramachandran outliers	78287	1008 (3.48-3.24)
Sidechain outliers	78261	1007 (3.48-3.24)
RSRZ outliers	66119	1141 (3.50-3.22)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	990	
1	B	990	
2	C	263	
2	E	263	
3	D	239	
3	F	239	

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 20939 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Insulin-degrading enzyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	932	Total	C	N	O	S	0	0	0
			7622	4911	1274	1415	22			
1	B	861	Total	C	N	O	S	0	0	0
			7040	4550	1173	1296	21			

There are 50 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	30	MET	-	EXPRESSION TAG	UNP P14735
A	31	HIS	-	EXPRESSION TAG	UNP P14735
A	32	HIS	-	EXPRESSION TAG	UNP P14735
A	33	HIS	-	EXPRESSION TAG	UNP P14735
A	34	HIS	-	EXPRESSION TAG	UNP P14735
A	35	HIS	-	EXPRESSION TAG	UNP P14735
A	36	HIS	-	EXPRESSION TAG	UNP P14735
A	37	ALA	-	EXPRESSION TAG	UNP P14735
A	38	ALA	-	EXPRESSION TAG	UNP P14735
A	39	GLY	-	EXPRESSION TAG	UNP P14735
A	40	ILE	-	EXPRESSION TAG	UNP P14735
A	41	PRO	-	EXPRESSION TAG	UNP P14735
A	110	LEU	CYS	CONFLICT	UNP P14735
A	171	SER	CYS	CONFLICT	UNP P14735
A	178	ALA	CYS	CONFLICT	UNP P14735
A	257	VAL	CYS	CONFLICT	UNP P14735
A	414	LEU	CYS	CONFLICT	UNP P14735
A	573	ASN	CYS	CONFLICT	UNP P14735
A	590	SER	CYS	CONFLICT	UNP P14735
A	789	SER	CYS	CONFLICT	UNP P14735
A	812	ALA	CYS	CONFLICT	UNP P14735
A	819	ALA	CYS	CONFLICT	UNP P14735
A	904	SER	CYS	CONFLICT	UNP P14735
A	966	ASN	CYS	CONFLICT	UNP P14735
A	974	ALA	CYS	CONFLICT	UNP P14735

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Chain	Residue	Modelled	Actual	Comment	Reference
B	30	MET	-	EXPRESSION TAG	UNP P14735
B	31	HIS	-	EXPRESSION TAG	UNP P14735
B	32	HIS	-	EXPRESSION TAG	UNP P14735
B	33	HIS	-	EXPRESSION TAG	UNP P14735
B	34	HIS	-	EXPRESSION TAG	UNP P14735
B	35	HIS	-	EXPRESSION TAG	UNP P14735
B	36	HIS	-	EXPRESSION TAG	UNP P14735
B	37	ALA	-	EXPRESSION TAG	UNP P14735
B	38	ALA	-	EXPRESSION TAG	UNP P14735
B	39	GLY	-	EXPRESSION TAG	UNP P14735
B	40	ILE	-	EXPRESSION TAG	UNP P14735
B	41	PRO	-	EXPRESSION TAG	UNP P14735
B	110	LEU	CYS	CONFLICT	UNP P14735
B	171	SER	CYS	CONFLICT	UNP P14735
B	178	ALA	CYS	CONFLICT	UNP P14735
B	257	VAL	CYS	CONFLICT	UNP P14735
B	414	LEU	CYS	CONFLICT	UNP P14735
B	573	ASN	CYS	CONFLICT	UNP P14735
B	590	SER	CYS	CONFLICT	UNP P14735
B	789	SER	CYS	CONFLICT	UNP P14735
B	812	ALA	CYS	CONFLICT	UNP P14735
B	819	ALA	CYS	CONFLICT	UNP P14735
B	904	SER	CYS	CONFLICT	UNP P14735
B	966	ASN	CYS	CONFLICT	UNP P14735
B	974	ALA	CYS	CONFLICT	UNP P14735

- Molecule 2 is a protein called Fab-bound IDE, heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	216	Total	C	N	O	S	0	0	0
			1637	1043	268	319	7			
2	E	211	Total	C	N	O	S	0	0	0
			1602	1023	263	309	7			

- Molecule 3 is a protein called Fab-bound IDE, light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	D	199	Total	C	N	O	S	0	0	0
			1528	962	255	306	5			
3	F	197	Total	C	N	O	S	0	0	0
			1508	947	250	306	5			

- Molecule 4 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	Zn 1	0	0

- Molecule 5 is water.

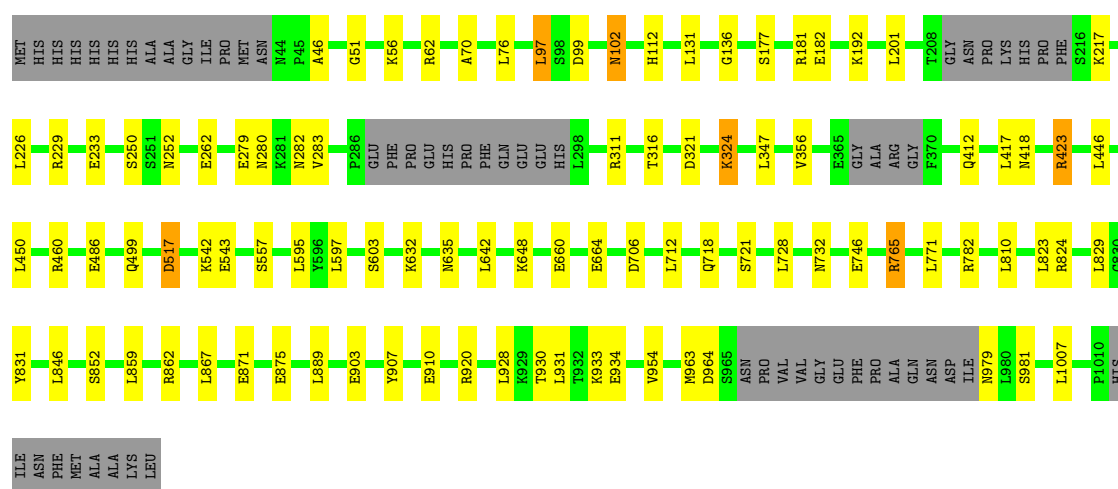
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total 1	O 1	0	0

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

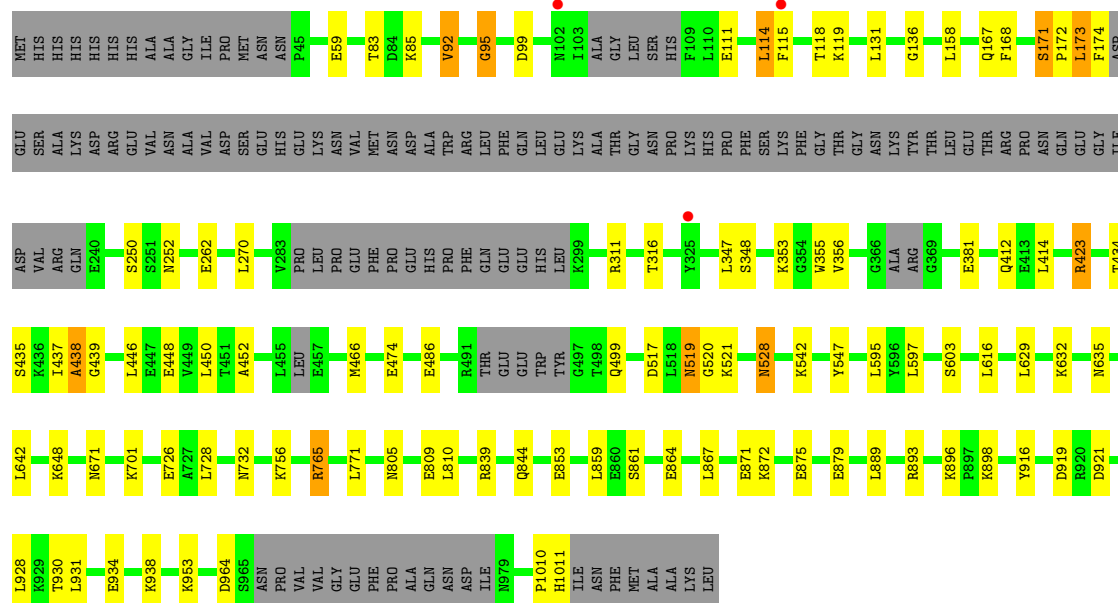
• Molecule 1: Insulin-degrading enzyme

Chain A:



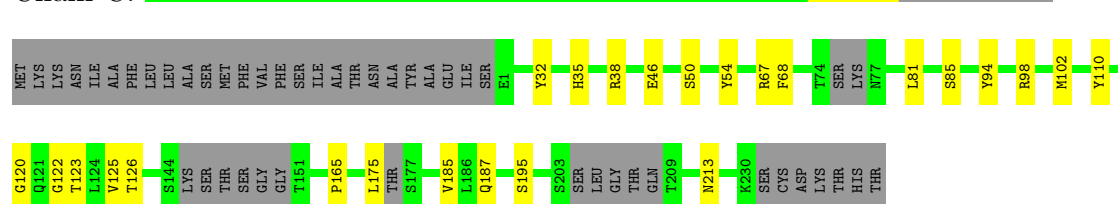
• Molecule 1: Insulin-degrading enzyme

Chain B:



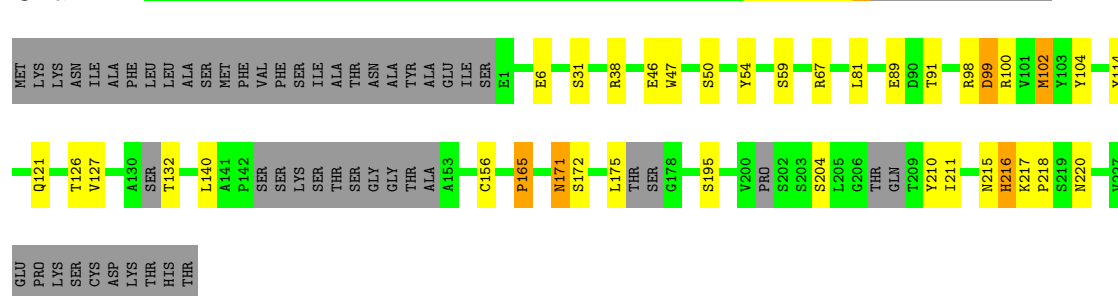
- Molecule 2: Fab-bound IDE, heavy chain

Chain C:



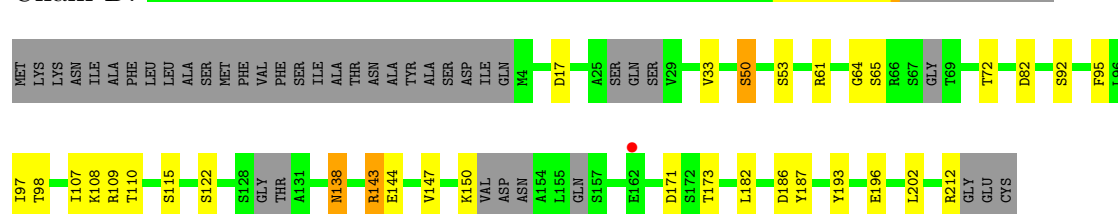
- Molecule 2: Fab-bound IDE, heavy chain

Chain E:



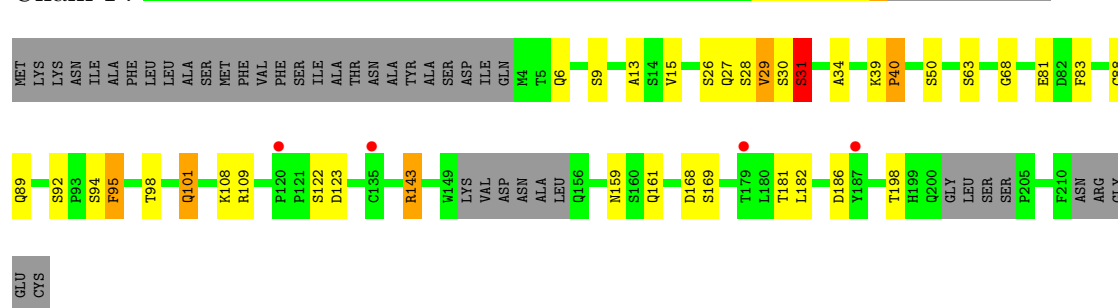
- Molecule 3: Fab-bound IDE, light chain

Chain D:



- Molecule 3: Fab-bound IDE, light chain

Chain F:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	56.95Å 131.66Å 377.04Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	49.85 – 3.35 49.85 – 3.35	Depositor EDS
% Data completeness (in resolution range)	99.9 (49.85-3.35) 95.1 (49.85-3.35)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	0.12	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.88 (at 3.33Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.1_1168)	Depositor
R, R_{free}	0.232 , 0.280 0.234 , 0.277	Depositor DCC
R_{free} test set	1988 reflections (4.99%)	DCC
Wilson B-factor (Å ²)	45.9	Xtriage
Anisotropy	0.237	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.24 , -10.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Outliers	0 of 40252 reflections	Xtriage
F_o, F_c correlation	0.83	EDS
Total number of atoms	20939	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.83% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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

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	0.32	0/7803	0.68	4/10549 (0.0%)
1	B	0.34	0/7206	0.72	6/9733 (0.1%)
2	C	0.34	0/1679	0.88	1/2285 (0.0%)
2	E	0.50	0/1641	0.85	2/2229 (0.1%)
3	D	0.33	0/1557	0.83	3/2106 (0.1%)
3	F	0.40	0/1540	0.90	5/2088 (0.2%)
All	All	0.35	0/21426	0.75	21/28990 (0.1%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	280	ASN	CB-CA-C	-15.27	79.86	110.40
1	B	171	SER	C-N-CD	-13.15	91.66	120.60
3	F	29	VAL	CB-CA-C	-11.61	89.35	111.40
2	E	220	ASN	CB-CA-C	-9.30	91.81	110.40
1	A	51	GLY	N-CA-C	-8.68	91.40	113.10
1	B	92	VAL	CB-CA-C	-7.74	96.69	111.40
1	A	97	LEU	CA-CB-CG	6.74	130.80	115.30
1	B	438	ALA	CB-CA-C	-6.40	100.50	110.10
3	F	31	SER	N-CA-C	-6.24	94.17	111.00
1	B	114	LEU	CB-CA-C	-6.11	98.59	110.20
3	D	64	GLY	N-CA-C	-6.09	97.86	113.10
3	D	50	SER	N-CA-C	-5.98	94.84	111.00
3	F	101	GLN	N-CA-C	-5.96	94.90	111.00
1	A	279	GLU	CB-CA-C	-5.92	98.56	110.40
3	F	50	SER	N-CA-C	-5.85	95.21	111.00
2	E	114	TYR	N-CA-C	-5.72	95.56	111.00
2	C	120	GLY	N-CA-C	-5.43	99.52	113.10
3	F	182	LEU	CA-CB-CG	5.39	127.70	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	439	GLY	N-CA-C	-5.34	99.75	113.10
1	B	95	GLY	N-CA-C	5.21	126.13	113.10
3	D	92	SER	CB-CA-C	5.15	119.89	110.10

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7622	0	0	28	0
1	B	7040	0	3	39	0
2	C	1637	0	0	6	0
2	E	1602	0	0	14	0
3	D	1528	0	0	5	0
3	F	1508	0	0	16	0
4	A	1	0	0	0	0
5	B	1	0	0	0	0
All	All	20939	0	3	105	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 5.

All (105) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
2:E:99:ASP:N	2:E:99:ASP:OD1	2.19	0.75
1:A:460:ARG:NH2	2:C:102:MET:O	2.20	0.74
1:B:115:PHE:CD1	1:B:168:PHE:CE2	2.76	0.73
1:B:809:GLU:OE2	1:B:839:ARG:NH2	2.23	0.70
1:B:893:ARG:NH2	1:B:921:ASP:OD1	2.25	0.70
2:E:216:HIS:ND1	2:E:218:PRO:CD	2.54	0.70
1:A:229:ARG:NE	1:A:233:GLU:OE2	2.25	0.69
1:A:517:ASP:N	1:A:517:ASP:OD1	2.25	0.68
3:F:159:ASN:ND2	3:F:181:THR:O	2.26	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:671:ASN:OD1	1:B:701:LYS:NZ	2.29	0.66
1:B:118:THR:CG2	1:B:171:SER:O	2.46	0.63
2:C:67:ARG:O	2:C:68:PHE:CD1	2.50	0.63
1:B:114:LEU:O	1:B:114:LEU:CG	2.47	0.63
1:A:99:ASP:O	1:A:217:LYS:NZ	2.32	0.62
1:B:118:THR:CG2	1:B:167:GLN:CB	2.78	0.61
2:E:38:ARG:NE	2:E:46:GLU:OE1	2.33	0.61
1:A:910:GLU:OE1	1:A:920:ARG:NH1	2.35	0.59
1:A:557:SER:OG	1:A:746:GLU:OE2	2.20	0.59
1:B:635:ASN:ND2	1:B:732:ASN:O	2.36	0.59
2:E:216:HIS:CE1	2:E:218:PRO:CD	2.86	0.58
1:A:782:ARG:NH1	1:A:963:MET:O	2.36	0.57
1:A:862:ARG:NH2	1:A:981:SER:O	2.38	0.57
1:A:706:ASP:OD1	1:B:756:LYS:NZ	2.38	0.56
2:E:91:THR:OG1	2:E:127:VAL:N	2.38	0.56
3:F:6:GLN:NE2	3:F:88:CYS:SG	2.78	0.56
1:B:99:ASP:OD1	1:B:99:ASP:N	2.38	0.56
1:A:423:ARG:NH1	1:A:423:ARG:CG	2.67	0.56
3:D:143:ARG:NE	3:D:143:ARG:O	2.38	0.56
1:A:321:ASP:O	1:A:324:LYS:NZ	2.39	0.55
1:A:311:ARG:NH2	1:A:664:GLU:OE2	2.39	0.55
2:C:67:ARG:C	2:C:68:PHE:CD1	2.81	0.54
3:D:50:SER:OG	3:D:53:SER:OG	2.24	0.54
3:F:29:VAL:CG1	3:F:29:VAL:O	2.54	0.54
1:B:864:GLU:OE1	1:B:953:LYS:NZ	2.41	0.54
1:A:635:ASN:ND2	1:A:732:ASN:O	2.41	0.54
2:C:38:ARG:NE	2:C:46:GLU:OE1	2.41	0.53
1:B:896:LYS:O	1:B:898:LYS:NZ	2.41	0.53
1:B:726:GLU:OE2	1:B:916:TYR:OH	2.27	0.52
3:F:27:GLN:CG	3:F:28:SER:N	2.73	0.52
1:B:131:LEU:O	1:B:136:GLY:N	2.43	0.52
1:B:434:THR:O	1:B:437:ILE:O	2.28	0.52
3:F:39:LYS:NZ	3:F:81:GLU:O	2.43	0.52
1:A:963:MET:SD	1:A:964:ASP:N	2.84	0.51
1:B:519:ASN:C	1:B:519:ASN:ND2	2.64	0.51
1:A:875:GLU:O	1:A:933:LYS:NZ	2.43	0.51
1:A:903:GLU:OE1	1:A:907:TYR:OH	2.29	0.51
1:B:423:ARG:NH1	1:B:423:ARG:CG	2.73	0.51
2:C:35:HIS:ND1	2:C:50:SER:OG	2.44	0.51
2:E:204:SER:OG	2:E:210:TYR:OH	2.29	0.50
1:B:92:VAL:CG1	1:B:92:VAL:O	2.59	0.50
3:F:13:ALA:O	3:F:108:LYS:N	2.44	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:102:ASN:O	1:A:102:ASN:ND2	2.45	0.50
1:A:177:SER:OG	1:A:181:ARG:NH1	2.45	0.50
1:B:311:ARG:NH2	1:B:381:GLU:OE1	2.45	0.50
1:B:250:SER:OG	1:B:252:ASN:OD1	2.30	0.49
3:D:61:ARG:NE	3:D:82:ASP:OD2	2.46	0.49
2:E:104:TYR:CD1	2:E:104:TYR:C	2.85	0.49
2:C:32:TYR:CZ	2:C:98:ARG:NH2	2.82	0.48
1:A:718:GLN:O	1:A:721:SER:OG	2.32	0.47
2:E:216:HIS:CE1	2:E:218:PRO:CG	2.97	0.47
2:E:171:ASN:OD1	2:E:211:ILE:N	2.47	0.47
1:A:56:LYS:NZ	1:A:62:ARG:O	2.47	0.47
1:A:829:LEU:O	1:A:852:SER:OG	2.33	0.47
1:A:46:ALA:O	1:A:70:ALA:N	2.47	0.46
1:A:250:SER:OG	1:A:252:ASN:OD1	2.33	0.46
1:B:59:GLU:O	1:B:59:GLU:CD	2.53	0.46
1:B:528:ASN:ND2	1:B:528:ASN:C	2.69	0.46
1:B:115:PHE:CD1	1:B:168:PHE:CD2	3.04	0.45
3:D:115:SER:CB	3:D:138:ASN:CB	2.94	0.45
1:B:173:LEU:CD1	1:B:174:PHE:C	2.86	0.45
1:A:765:ARG:NH1	1:A:765:ARG:CG	2.80	0.45
1:B:519:ASN:ND2	1:B:521:LYS:N	2.65	0.45
1:B:765:ARG:CG	1:B:765:ARG:NH1	2.80	0.44
2:E:217:LYS:N	2:E:218:PRO:CD	2.81	0.44
3:F:29:VAL:O	3:F:31:SER:N	2.50	0.44
1:B:173:LEU:O	1:B:174:PHE:CD1	2.70	0.44
2:E:47:TRP:NE1	2:E:50:SER:OG	2.51	0.44
2:E:31:SER:O	2:E:102:MET:O	2.36	0.44
2:E:6:GLU:OE1	2:E:6:GLU:N	2.51	0.44
1:B:448:GLU:O	1:B:452:ALA:N	2.50	0.44
1:B:83:THR:OG1	1:B:85:LYS:O	2.36	0.43
1:A:112:HIS:ND1	1:A:182:GLU:OE1	2.51	0.43
1:A:262:GLU:OE2	1:A:262:GLU:N	2.52	0.43
3:F:168:ASP:OD1	3:F:169:SER:N	2.52	0.43
1:B:547:TYR:OH	1:B:919:ASP:OD1	2.37	0.43
1:B:115:PHE:CE1	1:B:168:PHE:CE2	3.07	0.43
3:F:143:ARG:CB	3:F:143:ARG:NH1	2.82	0.42
1:B:805:ASN:OD1	1:B:844:GLN:NE2	2.53	0.42
1:B:173:LEU:CD2	1:B:173:LEU:C	2.85	0.42
1:B:353:LYS:NZ	1:B:355:TRP:CZ2	2.88	0.42
1:A:824:ARG:NH1	1:A:831:TYR:O	2.53	0.42
3:F:39:LYS:NZ	3:F:83:PHE:CE1	2.88	0.42
2:E:59:SER:OG	3:F:95:PHE:O	2.38	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
3:F:30:SER:OG	3:F:68:GLY:O	2.37	0.42
1:B:118:THR:CG2	1:B:119:LYS:N	2.81	0.41
3:F:29:VAL:C	3:F:31:SER:N	2.72	0.41
1:B:437:ILE:O	1:B:438:ALA:CB	2.69	0.41
3:F:159:ASN:ND2	3:F:181:THR:OG1	2.53	0.41
3:F:83:PHE:CG	3:F:83:PHE:O	2.73	0.41
3:D:187:TYR:O	3:D:193:TYR:OH	2.39	0.41
1:B:519:ASN:ND2	1:B:520:GLY:N	2.69	0.41
1:A:131:LEU:O	1:A:136:GLY:N	2.54	0.41
3:F:34:ALA:O	3:F:89:GLN:N	2.55	0.40
1:B:474:GLU:OE2	1:B:474:GLU:N	2.54	0.40
1:B:262:GLU:N	1:B:262:GLU:OE2	2.54	0.40

There are no symmetry-related clashes.

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	922/990 (93%)	886 (96%)	35 (4%)	1 (0%)	59	95
1	B	845/990 (85%)	804 (95%)	38 (4%)	3 (0%)	43	89
2	C	206/263 (78%)	188 (91%)	15 (7%)	3 (2%)	15	64
2	E	199/263 (76%)	178 (89%)	19 (10%)	2 (1%)	22	74
3	D	187/239 (78%)	165 (88%)	21 (11%)	1 (0%)	38	85
3	F	191/239 (80%)	162 (85%)	28 (15%)	1 (0%)	38	85
All	All	2550/2984 (86%)	2383 (94%)	156 (6%)	11 (0%)	43	89

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	D	138	ASN
1	A	282	ASN
1	B	95	GLY

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Mol	Chain	Res	Type
1	B	172	PRO
1	B	1010	PRO
2	C	122	GLY
3	F	40	PRO
2	C	165	PRO
2	E	165	PRO
2	E	171	ASN
2	C	85	SER

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	832/879 (95%)	785 (94%)	47 (6%)	30	73
1	B	768/879 (87%)	719 (94%)	49 (6%)	25	69
2	C	181/220 (82%)	169 (93%)	12 (7%)	24	68
2	E	176/220 (80%)	157 (89%)	19 (11%)	9	38
3	D	178/210 (85%)	155 (87%)	23 (13%)	6	29
3	F	176/210 (84%)	158 (90%)	18 (10%)	11	42
All	All	2311/2618 (88%)	2143 (93%)	168 (7%)	20	63

All (168) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	76	LEU
1	A	97	LEU
1	A	102	ASN
1	A	192	LYS
1	A	201	LEU
1	A	226	LEU
1	A	283	VAL
1	A	316	THR
1	A	324	LYS
1	A	347	LEU
1	A	356	VAL

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Mol	Chain	Res	Type
1	A	412	GLN
1	A	417	LEU
1	A	418	ASN
1	A	423	ARG
1	A	446	LEU
1	A	450	LEU
1	A	486	GLU
1	A	499	GLN
1	A	517	ASP
1	A	542	LYS
1	A	543	GLU
1	A	595	LEU
1	A	597	LEU
1	A	603	SER
1	A	632	LYS
1	A	642	LEU
1	A	648	LYS
1	A	660	GLU
1	A	712	LEU
1	A	728	LEU
1	A	765	ARG
1	A	771	LEU
1	A	810	LEU
1	A	823	LEU
1	A	846	LEU
1	A	859	LEU
1	A	867	LEU
1	A	871	GLU
1	A	889	LEU
1	A	928	LEU
1	A	930	THR
1	A	931	LEU
1	A	934	GLU
1	A	954	VAL
1	A	979	ASN
1	A	1007	LEU
1	B	111	GLU
1	B	158	LEU
1	B	173	LEU
1	B	270	LEU
1	B	316	THR
1	B	347	LEU

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Mol	Chain	Res	Type
1	B	348	SER
1	B	356	VAL
1	B	412	GLN
1	B	414	LEU
1	B	423	ARG
1	B	435	SER
1	B	446	LEU
1	B	450	LEU
1	B	466	MET
1	B	486	GLU
1	B	499	GLN
1	B	517	ASP
1	B	519	ASN
1	B	528	ASN
1	B	542	LYS
1	B	595	LEU
1	B	597	LEU
1	B	603	SER
1	B	616	LEU
1	B	629	LEU
1	B	632	LYS
1	B	642	LEU
1	B	648	LYS
1	B	728	LEU
1	B	765	ARG
1	B	771	LEU
1	B	810	LEU
1	B	853	GLU
1	B	859	LEU
1	B	861	SER
1	B	867	LEU
1	B	871	GLU
1	B	872	LYS
1	B	875	GLU
1	B	879	GLU
1	B	889	LEU
1	B	928	LEU
1	B	930	THR
1	B	931	LEU
1	B	934	GLU
1	B	938	LYS
1	B	964	ASP

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Mol	Chain	Res	Type
1	B	1011	HIS
2	C	54	TYR
2	C	81	LEU
2	C	94	TYR
2	C	110	TYR
2	C	123	THR
2	C	125	VAL
2	C	126	THR
2	C	175	LEU
2	C	185	VAL
2	C	187	GLN
2	C	195	SER
2	C	213	ASN
3	D	17	ASP
3	D	33	VAL
3	D	65	SER
3	D	72	THR
3	D	95	PHE
3	D	97	ILE
3	D	98	THR
3	D	107	ILE
3	D	108	LYS
3	D	109	ARG
3	D	110	THR
3	D	122	SER
3	D	143	ARG
3	D	144	GLU
3	D	147	VAL
3	D	150	LYS
3	D	171	ASP
3	D	173	THR
3	D	182	LEU
3	D	186	ASP
3	D	196	GLU
3	D	202	LEU
3	D	212	ARG
2	E	54	TYR
2	E	67	ARG
2	E	81	LEU
2	E	89	GLU
2	E	98	ARG
2	E	99	ASP

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Mol	Chain	Res	Type
2	E	100	ARG
2	E	102	MET
2	E	121	GLN
2	E	126	THR
2	E	132	THR
2	E	140	LEU
2	E	156	CYS
2	E	165	PRO
2	E	172	SER
2	E	175	LEU
2	E	195	SER
2	E	215	ASN
2	E	216	HIS
3	F	9	SER
3	F	15	VAL
3	F	26	SER
3	F	31	SER
3	F	40	PRO
3	F	63	SER
3	F	92	SER
3	F	94	SER
3	F	95	PHE
3	F	98	THR
3	F	101	GLN
3	F	109	ARG
3	F	122	SER
3	F	123	ASP
3	F	143	ARG
3	F	161	GLN
3	F	186	ASP
3	F	198	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

5.7 Other polymers ⓘ

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 ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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	932/990 (94%)	-0.06	0 100 100	12, 39, 101, 152	0
1	B	861/990 (86%)	-0.02	3 (0%) 91 72	11, 41, 100, 151	0
2	C	216/263 (82%)	0.07	0 100 100	19, 56, 97, 112	0
2	E	211/263 (80%)	0.03	0 100 100	18, 55, 98, 144	0
3	D	199/239 (83%)	0.19	1 (0%) 88 59	22, 55, 102, 138	0
3	F	197/239 (82%)	0.26	4 (2%) 62 25	20, 69, 116, 147	0
All	All	2616/2984 (87%)	0.01	8 (0%) 91 72	11, 46, 103, 152	0

All (8) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
3	F	179	THR	3.0
3	F	135	CYS	2.9
3	F	187	TYR	2.4
1	B	102	ASN	2.3
3	D	162	GLU	2.2
3	F	120	PRO	2.1
1	B	115	PHE	2.1
1	B	325	TYR	2.1

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	ZN	A	1101	1/1	0.06	-3.24	75,75,75,75	0

6.5 Other polymers

There are no such residues in this entry.