



Full wwPDB X-ray Structure Validation Report ⓘ

Aug 14, 2023 – 09:52 PM EDT

PDB ID : 1RY7
Title : Crystal Structure of the 3 Ig form of FGFR3c in complex with FGF1
Authors : Olsen, S.K.; Ibrahimi, O.A.; Raucci, A.; Zhang, F.; Eliseenkova, A.V.; Yayon, A.; Basilico, C.; Linhardt, R.J.; Schlessinger, J.; Mohammadi, M.
Deposited on : 2003-12-19
Resolution : 3.20 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 : 4.02b-467
Xtriage (Phenix) : 1.13
EDS : 2.35
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35

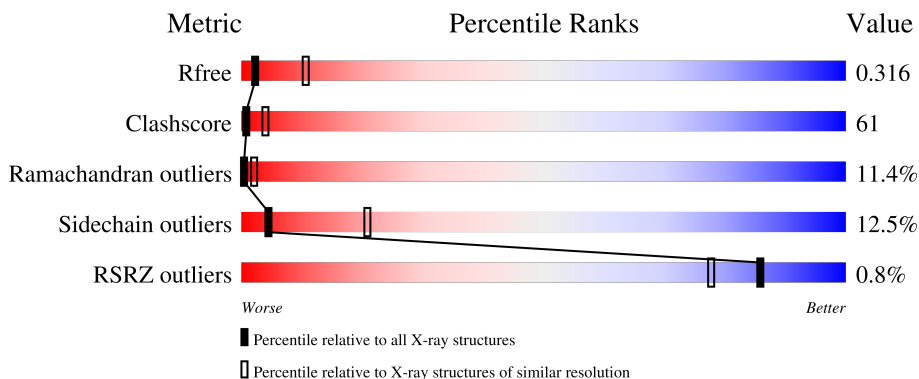
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.20 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}	130704	1133 (3.20-3.20)
Clashscore	141614	1253 (3.20-3.20)
Ramachandran outliers	138981	1234 (3.20-3.20)
Sidechain outliers	138945	1233 (3.20-3.20)
RSRZ outliers	127900	1095 (3.20-3.20)

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	155	 24% 61% 12% . .
2	B	334	 17% 35% 11% 36%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2766 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 Heparin-binding growth factor 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	151	1181	750	201	226	4	0	0	0

- Molecule 2 is a protein called Fibroblast growth factor receptor 3.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	B	213	1585	1010	277	292	6	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	32	MET	-	cloning artifact	UNP P22607

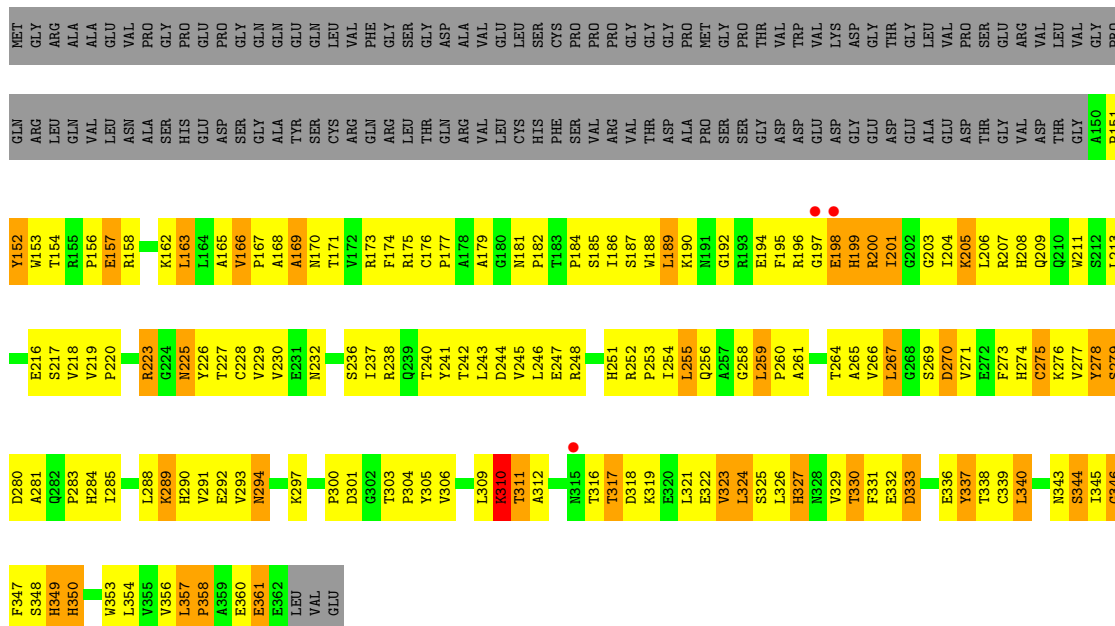
3 Residue-property plots

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: Heparin-binding growth factor 1



- Molecule 2: Fibroblast growth factor receptor 3



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	104.22Å 64.75Å 99.93Å 90.00° 94.67° 90.00°	Depositor
Resolution (Å)	25.00 – 3.20 29.07 – 3.12	Depositor EDS
% Data completeness (in resolution range)	(Not available) (25.00-3.20) 90.5 (29.07-3.12)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.17 (at 3.11Å)	Xtrriage
Refinement program	CNS	Depositor
R, R_{free}	0.277 , 0.341 0.258 , 0.316	Depositor DCC
R_{free} test set	1085 reflections (9.95%)	wwPDB-VP
Wilson B-factor (Å ²)	61.1	Xtrriage
Anisotropy	0.561	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 58.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.31$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	2766	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.42% 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](#)

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	0.40	0/1209	0.69	0/1638
2	B	0.55	2/1629 (0.1%)	0.75	2/2233 (0.1%)
All	All	0.49	2/2838 (0.1%)	0.73	2/3871 (0.1%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	170	ASN	CG-OD1	13.30	1.53	1.24
2	B	275	CYS	CB-SG	-5.32	1.73	1.81

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	170	ASN	CB-CG-OD1	-5.60	110.41	121.60
2	B	278	TYR	N-CA-C	-5.49	96.18	111.00

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	1181	0	1145	143	0
2	B	1585	0	1492	206	1
All	All	2766	0	2637	329	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 61.

All (329) 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)
2:B:255:LEU:HD22	2:B:277:VAL:HG22	1.36	1.04
1:A:36:HIS:HB3	1:A:49:THR:H	1.20	1.03
2:B:190:LYS:HB2	2:B:195:PHE:HD2	1.25	1.01
2:B:189:LEU:HD11	2:B:227:THR:HB	1.46	0.96
2:B:288:LEU:HD13	2:B:306:VAL:HG21	1.48	0.96
1:A:58:GLN:H	1:A:58:GLN:HE21	1.10	0.95
2:B:312:ALA:HB1	2:B:325:SER:O	1.71	0.92
2:B:297:LYS:HE2	2:B:304:PRO:HB3	1.52	0.90
2:B:173:ARG:HH11	2:B:173:ARG:HB3	1.39	0.88
2:B:310:LYS:O	2:B:311:THR:HG23	1.72	0.88
2:B:256:GLN:OE1	2:B:276:LYS:HD2	1.74	0.87
1:A:27:LYS:HD2	1:A:149:PRO:HB3	1.55	0.87
1:A:72:LYS:HD2	1:A:79:TYR:HE2	1.40	0.84
1:A:6:ILE:HG22	1:A:7:THR:H	1.40	0.83
2:B:225:ASN:HA	2:B:242:THR:HA	1.62	0.81
1:A:58:GLN:HE21	1:A:58:GLN:N	1.77	0.80
1:A:79:TYR:CE1	1:A:94:PRO:HB3	2.17	0.79
2:B:270:ASP:HB2	2:B:327:HIS:ND1	1.98	0.78
2:B:171:THR:HG23	2:B:216:GLU:HG2	1.65	0.77
1:A:39:ARG:HD3	1:A:54:ASP:OD2	1.85	0.76
1:A:85:ASP:HB2	1:A:87:LEU:HG	1.67	0.75
1:A:58:GLN:H	1:A:58:GLN:NE2	1.85	0.75
2:B:190:LYS:HB2	2:B:195:PHE:CD2	2.16	0.74
1:A:89:TYR:HD1	1:A:90:GLY:H	1.36	0.73
2:B:252:ARG:HG2	2:B:252:ARG:HH11	1.54	0.73
2:B:297:LYS:O	2:B:297:LYS:HD3	1.87	0.73
1:A:59:LEU:HA	1:A:72:LYS:O	1.89	0.73
2:B:256:GLN:NE2	2:B:259:LEU:HD22	2.03	0.72
1:A:34:GLY:HA3	1:A:36:HIS:CE1	2.25	0.72
1:A:81:ALA:HB2	1:A:98:CYS:HB3	1.71	0.72
1:A:36:HIS:HB3	1:A:49:THR:N	2.01	0.71
2:B:253:PRO:HD3	2:B:343:ASN:CB	2.20	0.71
2:B:253:PRO:HD3	2:B:343:ASN:HB3	1.73	0.71
2:B:176:CYS:HG	2:B:228:CYS:CB	2.04	0.71
1:A:41:LEU:HD12	1:A:45:THR:OG1	1.90	0.71
1:A:58:GLN:N	1:A:58:GLN:NE2	2.38	0.70
2:B:226:TYR:N	2:B:241:TYR:O	2.25	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:326:LEU:HD11	2:B:337:TYR:CE2	2.26	0.70
1:A:95:ASN:HD21	1:A:97:GLU:HB2	1.54	0.69
2:B:190:LYS:C	2:B:192:GLY:H	1.96	0.69
2:B:266:VAL:HG12	2:B:267:LEU:HD12	1.73	0.69
1:A:148:LEU:HD11	2:B:246:LEU:HD22	1.75	0.69
2:B:357:LEU:HD12	2:B:357:LEU:H	1.57	0.68
2:B:281:ALA:HB3	2:B:343:ASN:HD21	1.59	0.68
2:B:267:LEU:HD23	2:B:331:PHE:HE1	1.59	0.68
1:A:65:SER:HA	2:B:284:HIS:ND1	2.08	0.68
2:B:190:LYS:HD2	2:B:195:PHE:CE2	2.29	0.68
2:B:258:GLY:O	2:B:261:ALA:HB2	1.94	0.68
1:A:80:LEU:HD22	1:A:100:PHE:CE2	2.29	0.67
1:A:6:ILE:HG22	1:A:7:THR:N	2.10	0.67
1:A:108:HIS:HD2	2:B:248:ARG:HB2	1.59	0.67
2:B:275:CYS:HG	2:B:339:CYS:HG	1.25	0.66
1:A:79:TYR:HE1	1:A:94:PRO:HB3	1.61	0.66
1:A:108:HIS:CD2	2:B:248:ARG:HB2	2.31	0.66
2:B:190:LYS:HD2	2:B:195:PHE:HE2	1.61	0.66
1:A:103:ARG:HG2	1:A:103:ARG:HH11	1.60	0.66
1:A:138:THR:HA	1:A:142:GLN:OE1	1.95	0.66
1:A:95:ASN:ND2	1:A:97:GLU:HB2	2.11	0.65
1:A:27:LYS:CD	1:A:149:PRO:HB3	2.24	0.65
1:A:42:PRO:HG3	1:A:56:HIS:NE2	2.12	0.65
1:A:63:ALA:HB2	1:A:69:VAL:HG12	1.78	0.65
1:A:16:PHE:CD1	2:B:256:GLN:HB2	2.32	0.65
1:A:80:LEU:HD23	1:A:81:ALA:N	2.11	0.65
2:B:197:GLY:HA2	2:B:204:ILE:CD1	2.26	0.65
1:A:33:ASN:HB2	1:A:144:ALA:HA	1.79	0.65
2:B:156:PRO:HG2	2:B:157:GLU:H	1.62	0.65
2:B:229:VAL:HG11	2:B:238:ARG:HH21	1.62	0.65
1:A:72:LYS:HD2	1:A:79:TYR:CE2	2.27	0.64
1:A:76:THR:OG1	1:A:78:GLN:HG3	1.97	0.64
2:B:243:LEU:HD13	2:B:244:ASP:N	2.12	0.64
2:B:331:PHE:HA	2:B:356:VAL:CG2	2.27	0.64
1:A:51:ASP:C	1:A:53:SER:H	2.01	0.63
2:B:317:THR:HG22	2:B:318:ASP:N	2.13	0.63
1:A:105:GLU:O	1:A:108:HIS:N	2.31	0.63
1:A:36:HIS:HA	1:A:49:THR:O	1.99	0.63
2:B:171:THR:HA	2:B:216:GLU:HA	1.81	0.62
2:B:289:LYS:HE2	2:B:290:HIS:N	2.14	0.62
1:A:80:LEU:HD23	1:A:80:LEU:C	2.20	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:LEU:HD13	1:A:147:PHE:HE1	1.65	0.61
2:B:166:VAL:HG22	2:B:167:PRO:HD2	1.81	0.61
2:B:196:ARG:N	2:B:199:HIS:HB2	2.14	0.61
2:B:267:LEU:HD23	2:B:331:PHE:CE1	2.35	0.61
1:A:106:GLU:C	1:A:108:HIS:H	2.04	0.61
2:B:195:PHE:HE1	2:B:204:ILE:HG21	1.64	0.61
2:B:271:VAL:HG22	2:B:326:LEU:HB2	1.82	0.61
2:B:357:LEU:HD12	2:B:357:LEU:N	2.15	0.61
2:B:319:LYS:O	2:B:319:LYS:HD3	2.01	0.61
1:A:23:TYR:CD2	2:B:283:PRO:HG3	2.36	0.61
2:B:169:ALA:HA	2:B:217:SER:HA	1.83	0.61
1:A:95:ASN:O	1:A:97:GLU:N	2.33	0.60
2:B:173:ARG:HB3	2:B:173:ARG:NH1	2.13	0.60
2:B:273:PHE:O	2:B:323:VAL:HA	2.01	0.60
2:B:201:ILE:C	2:B:203:GLY:H	2.05	0.60
2:B:323:VAL:CG2	2:B:324:LEU:N	2.64	0.60
2:B:252:ARG:HG3	2:B:348:SER:HB2	1.83	0.60
2:B:312:ALA:CB	2:B:325:SER:O	2.47	0.60
2:B:197:GLY:HA2	2:B:204:ILE:HD11	1.84	0.60
1:A:28:LEU:HD22	1:A:57:ILE:HG13	1.82	0.59
1:A:39:ARG:NE	1:A:41:LEU:HD21	2.18	0.59
1:A:129:ASN:HD21	1:A:131:SER:HB3	1.66	0.59
1:A:89:TYR:CD1	1:A:90:GLY:N	2.71	0.59
1:A:103:ARG:HG2	1:A:103:ARG:NH1	2.16	0.59
1:A:133:LYS:HD3	1:A:138:THR:HG22	1.84	0.59
1:A:59:LEU:HD22	1:A:71:ILE:HG22	1.83	0.58
2:B:357:LEU:HB3	2:B:358:PRO:HD2	1.85	0.58
2:B:279:SER:O	2:B:280:ASP:C	2.42	0.58
2:B:157:GLU:HG2	2:B:158:ARG:N	2.18	0.58
2:B:186:ILE:HG23	2:B:186:ILE:O	2.04	0.57
2:B:333:ASP:OD1	2:B:333:ASP:N	2.36	0.57
1:A:11:ALA:O	1:A:12:LEU:HB2	2.03	0.57
2:B:223:ARG:CB	2:B:244:ASP:HA	2.34	0.57
2:B:343:ASN:OD1	2:B:344:SER:N	2.38	0.57
1:A:33:ASN:HD21	1:A:128:LYS:N	2.03	0.57
1:A:83:ASP:OD1	1:A:87:LEU:HD12	2.05	0.57
2:B:290:HIS:HA	2:B:306:VAL:HG12	1.85	0.56
1:A:59:LEU:HD23	1:A:72:LYS:O	2.05	0.56
2:B:348:SER:O	2:B:349:HIS:HB3	2.05	0.56
2:B:289:LYS:O	2:B:306:VAL:HA	2.06	0.56
2:B:254:ILE:C	2:B:255:LEU:HD23	2.26	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:HIS:CB	1:A:49:THR:H	2.06	0.55
2:B:176:CYS:SG	2:B:228:CYS:SG	3.03	0.55
2:B:265:ALA:O	2:B:356:VAL:HA	2.06	0.55
1:A:145:ILE:HG13	1:A:146:LEU:HD23	1.87	0.55
2:B:154:THR:OG1	2:B:177:PRO:HB2	2.07	0.55
2:B:195:PHE:HA	2:B:199:HIS:ND1	2.21	0.55
2:B:173:ARG:HH11	2:B:173:ARG:CB	2.17	0.54
1:A:18:LEU:HB3	2:B:278:TYR:HB2	1.88	0.54
2:B:190:LYS:O	2:B:192:GLY:N	2.38	0.54
1:A:85:ASP:N	1:A:85:ASP:OD2	2.39	0.54
1:A:51:ASP:O	1:A:53:SER:N	2.39	0.54
2:B:230:VAL:O	2:B:236:SER:HA	2.08	0.54
2:B:196:ARG:C	2:B:198:GLU:H	2.11	0.54
1:A:38:LEU:HD13	1:A:147:PHE:CE1	2.42	0.53
1:A:9:PHE:O	1:A:11:ALA:N	2.41	0.53
2:B:229:VAL:CG1	2:B:238:ARG:HH21	2.21	0.53
1:A:88:LEU:HD13	1:A:124:VAL:HG13	1.89	0.53
2:B:276:LYS:HG2	2:B:321:LEU:CD2	2.38	0.53
2:B:266:VAL:HG13	2:B:358:PRO:O	2.08	0.53
2:B:288:LEU:HD11	2:B:340:LEU:HD12	1.90	0.53
2:B:310:LYS:O	2:B:311:THR:CG2	2.53	0.53
2:B:187:SER:CB	2:B:229:VAL:HB	2.39	0.53
2:B:223:ARG:HB2	2:B:244:ASP:HA	1.91	0.53
1:A:37:PHE:O	1:A:38:LEU:C	2.45	0.52
1:A:34:GLY:HA3	1:A:36:HIS:HE1	1.71	0.52
2:B:189:LEU:HA	2:B:195:PHE:N	2.24	0.52
1:A:109:TYR:HD2	1:A:146:LEU:HD13	1.74	0.52
2:B:357:LEU:CB	2:B:358:PRO:HD2	2.38	0.52
2:B:264:THR:HG22	2:B:265:ALA:N	2.25	0.52
1:A:39:ARG:NH2	1:A:47:ASP:OD2	2.33	0.52
1:A:110:ASN:OD1	2:B:248:ARG:NH1	2.40	0.52
2:B:281:ALA:HB3	2:B:343:ASN:ND2	2.24	0.52
2:B:218:VAL:HB	2:B:245:VAL:HG21	1.92	0.51
1:A:79:TYR:CD1	1:A:94:PRO:HB3	2.45	0.51
1:A:108:HIS:HB3	2:B:248:ARG:HG3	1.92	0.51
2:B:151:PRO:HA	2:B:179:ALA:O	2.11	0.51
2:B:207:ARG:C	2:B:209:GLN:H	2.14	0.51
2:B:204:ILE:HG13	2:B:205:LYS:N	2.26	0.51
2:B:190:LYS:C	2:B:192:GLY:N	2.64	0.51
2:B:318:ASP:HB3	2:B:321:LEU:HB2	1.93	0.51
1:A:42:PRO:HG3	1:A:56:HIS:CD2	2.46	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:GLU:HG2	1:A:115:LYS:HD2	1.92	0.51
2:B:220:PRO:HA	2:B:245:VAL:HB	1.92	0.51
2:B:166:VAL:HG22	2:B:167:PRO:CD	2.41	0.51
2:B:309:LEU:HD13	2:B:324:LEU:HD11	1.92	0.51
2:B:255:LEU:HD23	2:B:255:LEU:N	2.26	0.50
2:B:162:LYS:CB	2:B:241:TYR:HA	2.41	0.50
2:B:276:LYS:HG2	2:B:321:LEU:HD21	1.93	0.50
2:B:337:TYR:CD1	2:B:337:TYR:N	2.76	0.50
1:A:30:TYR:HB3	1:A:148:LEU:O	2.11	0.50
2:B:195:PHE:CE1	2:B:204:ILE:HG21	2.46	0.50
1:A:39:ARG:O	1:A:46:VAL:HA	2.11	0.50
2:B:198:GLU:C	2:B:200:ARG:H	2.15	0.50
2:B:223:ARG:HB2	2:B:223:ARG:NH1	2.27	0.50
2:B:230:VAL:HB	2:B:237:ILE:HG13	1.93	0.50
2:B:270:ASP:C	2:B:270:ASP:OD2	2.50	0.49
2:B:253:PRO:HD3	2:B:343:ASN:HB2	1.92	0.49
1:A:51:ASP:C	1:A:53:SER:N	2.66	0.49
2:B:266:VAL:O	2:B:269:SER:OG	2.31	0.49
2:B:309:LEU:HD22	2:B:310:LYS:HE3	1.95	0.49
2:B:252:ARG:HH11	2:B:252:ARG:CG	2.24	0.49
2:B:186:ILE:HD12	2:B:211:TRP:HA	1.93	0.49
1:A:81:ALA:O	1:A:88:LEU:HA	2.12	0.48
1:A:6:ILE:CG2	1:A:7:THR:H	2.20	0.48
2:B:243:LEU:HD11	2:B:245:VAL:HG23	1.95	0.48
2:B:278:TYR:CD2	2:B:279:SER:N	2.82	0.48
2:B:309:LEU:CD1	2:B:324:LEU:HD11	2.43	0.48
1:A:65:SER:HA	2:B:284:HIS:CE1	2.48	0.48
2:B:340:LEU:HD22	2:B:347:PHE:HD2	1.79	0.48
1:A:33:ASN:ND2	1:A:128:LYS:N	2.60	0.48
2:B:153:TRP:O	2:B:156:PRO:HD3	2.13	0.48
2:B:360:GLU:O	2:B:361:GLU:CB	2.61	0.48
2:B:271:VAL:HG21	2:B:354:LEU:CD1	2.44	0.48
2:B:278:TYR:O	2:B:279:SER:HB2	2.14	0.48
1:A:36:HIS:ND1	1:A:36:HIS:N	2.61	0.48
2:B:337:TYR:N	2:B:337:TYR:HD1	2.11	0.48
1:A:89:TYR:HD1	1:A:90:GLY:N	2.03	0.47
2:B:152:TYR:N	2:B:152:TYR:CD1	2.81	0.47
2:B:176:CYS:SG	2:B:228:CYS:CB	3.01	0.47
1:A:79:TYR:HD1	1:A:94:PRO:CA	2.27	0.47
1:A:81:ALA:HB1	1:A:97:GLU:O	2.15	0.47
2:B:189:LEU:HD12	2:B:189:LEU:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:ASN:HD21	2:B:248:ARG:HH22	1.61	0.47
1:A:39:ARG:HD3	1:A:54:ASP:CG	2.35	0.47
2:B:252:ARG:HD2	2:B:347:PHE:O	2.14	0.47
2:B:349:HIS:C	2:B:350:HIS:HD1	2.17	0.47
2:B:171:THR:CG2	2:B:216:GLU:HG2	2.40	0.47
2:B:198:GLU:O	2:B:200:ARG:N	2.48	0.47
2:B:330:THR:C	2:B:332:GLU:H	2.17	0.47
2:B:336:GLU:HG3	2:B:353:TRP:CE2	2.50	0.47
2:B:310:LYS:HD2	2:B:310:LYS:N	2.30	0.47
1:A:18:LEU:CB	2:B:278:TYR:HB2	2.45	0.46
1:A:81:ALA:HB3	1:A:89:TYR:CE1	2.50	0.46
1:A:114:SER:O	1:A:115:LYS:C	2.53	0.46
2:B:220:PRO:O	2:B:223:ARG:HG2	2.15	0.46
1:A:134:ARG:O	1:A:135:GLY:C	2.54	0.46
2:B:255:LEU:O	2:B:256:GLN:C	2.53	0.46
1:A:113:ILE:O	1:A:115:LYS:N	2.49	0.46
2:B:338:THR:HG23	2:B:350:HIS:O	2.15	0.46
2:B:343:ASN:OD1	2:B:343:ASN:C	2.54	0.46
2:B:285:ILE:HG13	2:B:322:GLU:OE2	2.15	0.46
1:A:6:ILE:O	1:A:7:THR:HB	2.16	0.46
1:A:24:LYS:HD2	2:B:319:LYS:O	2.16	0.46
1:A:88:LEU:CD1	1:A:124:VAL:HG13	2.45	0.46
1:A:124:VAL:HG12	1:A:124:VAL:O	2.16	0.46
2:B:207:ARG:C	2:B:209:GLN:N	2.70	0.46
2:B:174:PHE:O	2:B:213:LEU:N	2.49	0.46
2:B:243:LEU:HD13	2:B:243:LEU:C	2.36	0.46
2:B:274:HIS:O	2:B:275:CYS:SG	2.74	0.46
1:A:15:LYS:O	2:B:256:GLN:HG3	2.16	0.46
1:A:41:LEU:HD11	1:A:47:ASP:OD1	2.16	0.45
2:B:293:VAL:O	2:B:294:ASN:C	2.54	0.45
1:A:83:ASP:HB3	1:A:89:TYR:HE2	1.81	0.45
2:B:153:TRP:CE2	2:B:237:ILE:HD12	2.51	0.45
2:B:219:VAL:HG13	2:B:220:PRO:HD2	1.98	0.45
2:B:277:VAL:O	2:B:277:VAL:HG12	2.15	0.45
1:A:79:TYR:O	1:A:80:LEU:C	2.55	0.45
1:A:83:ASP:HB3	1:A:89:TYR:CE2	2.51	0.45
1:A:103:ARG:HH11	1:A:103:ARG:CG	2.28	0.45
2:B:266:VAL:HB	2:B:269:SER:OG	2.17	0.45
2:B:152:TYR:CE1	2:B:179:ALA:HB3	2.52	0.45
1:A:109:TYR:CD2	1:A:146:LEU:HD13	2.50	0.45
2:B:206:LEU:HD12	2:B:206:LEU:C	2.37	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:246:LEU:HD12	2:B:246:LEU:N	2.31	0.45
2:B:259:LEU:HA	2:B:260:PRO:C	2.37	0.45
2:B:266:VAL:HG12	2:B:267:LEU:N	2.31	0.45
1:A:29:LEU:HD23	1:A:29:LEU:HA	1.80	0.45
2:B:253:PRO:CG	2:B:343:ASN:HB2	2.47	0.45
2:B:316:THR:HB	2:B:317:THR:H	1.56	0.44
2:B:230:VAL:HB	2:B:237:ILE:CG1	2.47	0.44
1:A:32:SER:HB2	1:A:146:LEU:HB2	1.99	0.44
2:B:223:ARG:HB2	2:B:223:ARG:CZ	2.47	0.44
2:B:174:PHE:CD1	2:B:241:TYR:CD1	3.06	0.44
2:B:330:THR:C	2:B:332:GLU:N	2.71	0.44
1:A:148:LEU:HD12	1:A:148:LEU:C	2.38	0.44
1:A:39:ARG:O	1:A:46:VAL:HG13	2.17	0.44
2:B:331:PHE:HA	2:B:356:VAL:HG21	1.99	0.44
1:A:6:ILE:CG2	1:A:7:THR:N	2.81	0.44
2:B:189:LEU:HA	2:B:195:PHE:H	1.82	0.44
2:B:196:ARG:CA	2:B:199:HIS:HB2	2.47	0.44
2:B:169:ALA:HA	2:B:217:SER:CA	2.46	0.43
2:B:326:LEU:HD11	2:B:337:TYR:CZ	2.52	0.43
2:B:345:ILE:O	2:B:346:GLY:O	2.35	0.43
1:A:80:LEU:HD22	1:A:100:PHE:HE2	1.80	0.43
1:A:95:ASN:H	1:A:98:CYS:HG	1.64	0.43
1:A:117:HIS:O	1:A:118:ALA:C	2.56	0.43
1:A:41:LEU:O	1:A:44:GLY:N	2.51	0.43
2:B:188:TRP:CD1	2:B:206:LEU:HD22	2.53	0.43
2:B:198:GLU:C	2:B:200:ARG:N	2.71	0.43
1:A:24:LYS:HB3	2:B:319:LYS:NZ	2.33	0.43
1:A:86:GLY:O	1:A:135:GLY:N	2.47	0.43
2:B:176:CYS:HG	2:B:228:CYS:HB2	1.78	0.43
2:B:201:ILE:C	2:B:203:GLY:N	2.71	0.43
1:A:29:LEU:O	1:A:30:TYR:O	2.37	0.43
1:A:54:ASP:O	1:A:57:ILE:HG12	2.18	0.43
1:A:106:GLU:C	1:A:108:HIS:N	2.71	0.43
2:B:289:LYS:HZ3	2:B:291:VAL:N	2.17	0.43
1:A:76:THR:OG1	1:A:77:GLY:N	2.51	0.43
2:B:293:VAL:HB	2:B:305:TYR:CE1	2.54	0.43
2:B:207:ARG:O	2:B:209:GLN:N	2.52	0.42
1:A:61:LEU:N	1:A:61:LEU:HD12	2.34	0.42
1:A:90:GLY:O	1:A:91:SER:O	2.37	0.42
2:B:151:PRO:HB3	2:B:184:PRO:HG3	2.01	0.42
1:A:39:ARG:CD	1:A:54:ASP:OD2	2.62	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:GLY:O	1:A:137:ARG:N	2.52	0.42
1:A:8:THR:O	1:A:12:LEU:HB2	2.19	0.42
1:A:17:ASN:C	1:A:18:LEU:HD22	2.40	0.42
1:A:90:GLY:O	1:A:91:SER:C	2.58	0.42
2:B:252:ARG:HG2	2:B:252:ARG:NH1	2.29	0.42
1:A:39:ARG:HG3	1:A:56:HIS:HB2	2.02	0.42
2:B:185:SER:O	2:B:230:VAL:HA	2.20	0.42
2:B:290:HIS:HB3	2:B:297:LYS:HE3	2.01	0.42
2:B:323:VAL:HG23	2:B:324:LEU:N	2.34	0.42
1:A:72:LYS:HG2	1:A:73:SER:N	2.35	0.42
1:A:105:GLU:HG3	1:A:111:THR:HG23	2.01	0.42
1:A:110:ASN:ND2	2:B:248:ARG:HH22	2.18	0.42
1:A:150:LEU:HD11	2:B:165:ALA:HB2	2.01	0.42
2:B:181:ASN:HA	2:B:182:PRO:HA	1.83	0.42
2:B:156:PRO:HG2	2:B:157:GLU:N	2.32	0.41
2:B:276:LYS:CG	2:B:321:LEU:HD21	2.50	0.41
2:B:349:HIS:C	2:B:349:HIS:CD2	2.94	0.41
1:A:9:PHE:O	1:A:10:THR:C	2.58	0.41
1:A:95:ASN:C	1:A:97:GLU:N	2.72	0.41
2:B:169:ALA:CA	2:B:217:SER:HA	2.51	0.41
2:B:252:ARG:CG	2:B:252:ARG:NH1	2.84	0.41
1:A:30:TYR:CE2	2:B:163:LEU:HD12	2.55	0.41
1:A:80:LEU:C	1:A:80:LEU:CD2	2.88	0.41
2:B:270:ASP:N	2:B:327:HIS:O	2.54	0.41
2:B:274:HIS:CG	2:B:321:LEU:HD13	2.54	0.41
2:B:291:VAL:HG22	2:B:292:GLU:N	2.35	0.41
2:B:223:ARG:HA	2:B:243:LEU:CD1	2.51	0.41
1:A:43:ASP:CG	1:A:44:GLY:N	2.74	0.41
1:A:28:LEU:HD13	1:A:37:PHE:CD1	2.56	0.41
1:A:155:ASP:OD1	2:B:223:ARG:CD	2.68	0.41
1:A:54:ASP:C	1:A:56:HIS:H	2.24	0.41
1:A:79:TYR:HD1	1:A:94:PRO:HA	1.86	0.41
2:B:227:THR:OG1	2:B:240:THR:HG23	2.21	0.41
2:B:289:LYS:HE2	2:B:290:HIS:H	1.85	0.41
2:B:301:ASP:C	2:B:303:THR:H	2.24	0.41
2:B:169:ALA:C	2:B:217:SER:HA	2.40	0.41
1:A:31:CYS:SG	1:A:126:LEU:HB2	2.61	0.40
1:A:99:LEU:HA	1:A:99:LEU:HD23	1.80	0.40
2:B:267:LEU:HG	2:B:331:PHE:CZ	2.56	0.40
1:A:16:PHE:HB2	2:B:256:GLN:OE1	2.21	0.40
1:A:127:LYS:O	1:A:130:GLY:N	2.35	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:82:MET:HE3	1:A:88:LEU:HD21	2.02	0.40
2:B:223:ARG:HA	2:B:243:LEU:HD12	2.02	0.40
2:B:288:LEU:HB2	2:B:338:THR:HB	2.04	0.40
2:B:175:ARG:HA	2:B:211:TRP:O	2.21	0.40
1:A:139:HIS:O	1:A:142:GLN:HG2	2.21	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)
2:B:327:HIS:CE1	2:B:327:HIS:CE1[2_656]	2.06	0.14

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	149/155 (96%)	105 (70%)	23 (15%)	21 (14%)	0	1
2	B	211/334 (63%)	147 (70%)	44 (21%)	20 (10%)	0	3
All	All	360/489 (74%)	252 (70%)	67 (19%)	41 (11%)	0	2

All (41) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	12	LEU
1	A	30	TYR
1	A	51	ASP
1	A	52	ARG
1	A	80	LEU
1	A	96	GLU
1	A	114	SER
2	B	169	ALA

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Mol	Chain	Res	Type
2	B	198	GLU
2	B	205	LYS
2	B	311	THR
2	B	317	THR
2	B	358	PRO
2	B	361	GLU
1	A	7	THR
1	A	55	GLN
1	A	91	SER
1	A	108	HIS
2	B	194	GLU
2	B	199	HIS
2	B	279	SER
2	B	294	ASN
2	B	346	GLY
1	A	32	SER
1	A	77	GLY
1	A	84	THR
1	A	128	LYS
2	B	208	HIS
2	B	270	ASP
2	B	310	LYS
1	A	9	PHE
1	A	135	GLY
2	B	168	ALA
2	B	201	ILE
2	B	225	ASN
2	B	300	PRO
1	A	14	GLU
2	B	329	VAL
1	A	49	THR
1	A	136	PRO
1	A	35	GLY

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	128/135 (95%)	118 (92%)	10 (8%)	12	43
2	B	161/278 (58%)	135 (84%)	26 (16%)	2	11
All	All	289/413 (70%)	253 (88%)	36 (12%)	4	21

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	36	HIS
1	A	43	ASP
1	A	49	THR
1	A	58	GLN
1	A	64	GLU
1	A	66	VAL
1	A	84	THR
1	A	85	ASP
1	A	104	LEU
1	A	132	CYS
2	B	152	TYR
2	B	157	GLU
2	B	163	LEU
2	B	166	VAL
2	B	189	LEU
2	B	200	ARG
2	B	223	ARG
2	B	232	ASN
2	B	247	GLU
2	B	251	HIS
2	B	255	LEU
2	B	259	LEU
2	B	267	LEU
2	B	289	LYS
2	B	310	LYS
2	B	323	VAL
2	B	324	LEU
2	B	327	HIS
2	B	330	THR
2	B	333	ASP
2	B	337	TYR
2	B	340	LEU
2	B	344	SER
2	B	349	HIS
2	B	350	HIS

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Mol	Chain	Res	Type
2	B	357	LEU

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	33	ASN
1	A	58	GLN
1	A	95	ASN
1	A	129	ASN
2	B	232	ASN
2	B	286	GLN
2	B	294	ASN
2	B	349	HIS

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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	151/155 (97%)	-0.42	0 100 100	7, 39, 65, 71	0
2	B	213/334 (63%)	-0.13	3 (1%) 75 63	2, 46, 82, 91	0
All	All	364/489 (74%)	-0.25	3 (0%) 86 78	2, 44, 74, 91	0

All (3) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	197	GLY	4.2
2	B	198	GLU	2.5
2	B	315	ASN	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 monosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.