



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:13 PM UTC

PDB ID : 4GOV / pdb_00004gov
Title : The crystal structure of human fascin 1 S39D mutant
Authors : Yang, S.Y.; Huang, F.K.; Huang, J.; Chen, S.; Jakoncic, J.; Leo-Macias, A.;
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Deposited on : 2012-08-20
Resolution : 2.20 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	:	4-5-2 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

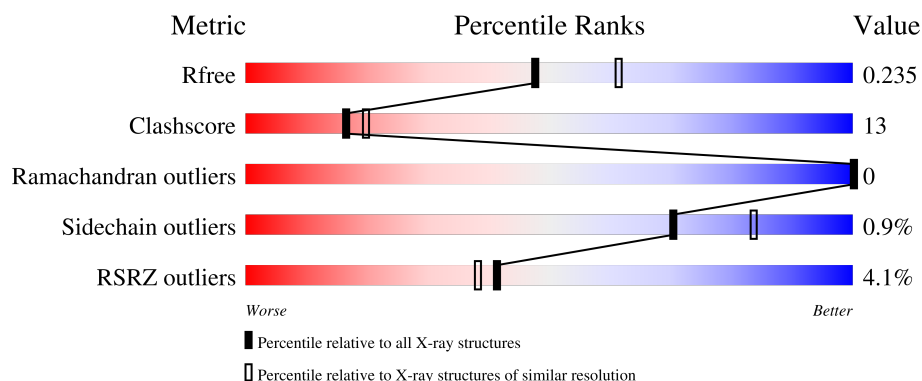
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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}	180053	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	493	6% 72% 22% • 5%
1	B	493	2% 72% 25% •

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
2	BR	A	502	-	-	X	-
4	GOL	B	518	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 8114 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 Fascin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	470	Total	C	N	O	S	0	1	0
			3675	2301	651	708	15			
1	B	483	Total	C	N	O	S	0	4	0
			3791	2371	676	727	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	39	ASP	SER	engineered mutation	UNP Q16658
B	39	ASP	SER	engineered mutation	UNP Q16658

- Molecule 2 is BROMIDE ION (CCD ID: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Br	0	0
			2	2		
2	B	4	Total	Br	0	0
			4	4		

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

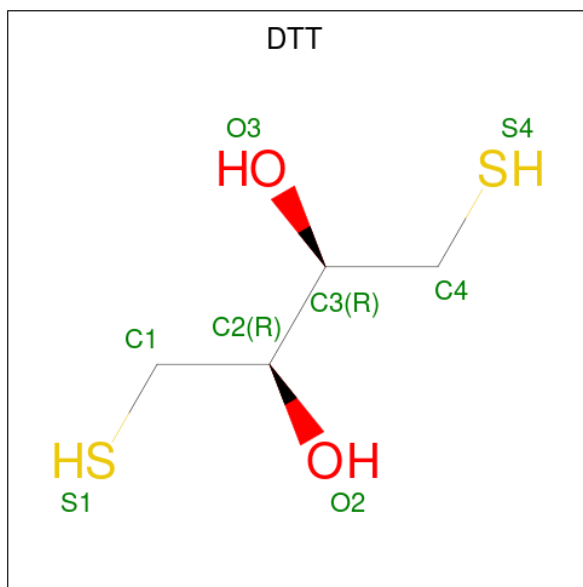
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	5	Total	Cl	0	0
			5	5		
3	B	8	Total	Cl	0	0
			8	8		

- Molecule 4 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is 2,3-DIHYDROXY-1,4-DITHIOBUTANE (CCD ID: DTT) (formula: $C_4H_{10}O_2S_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	O	S	0	0
			8	4	2	2		

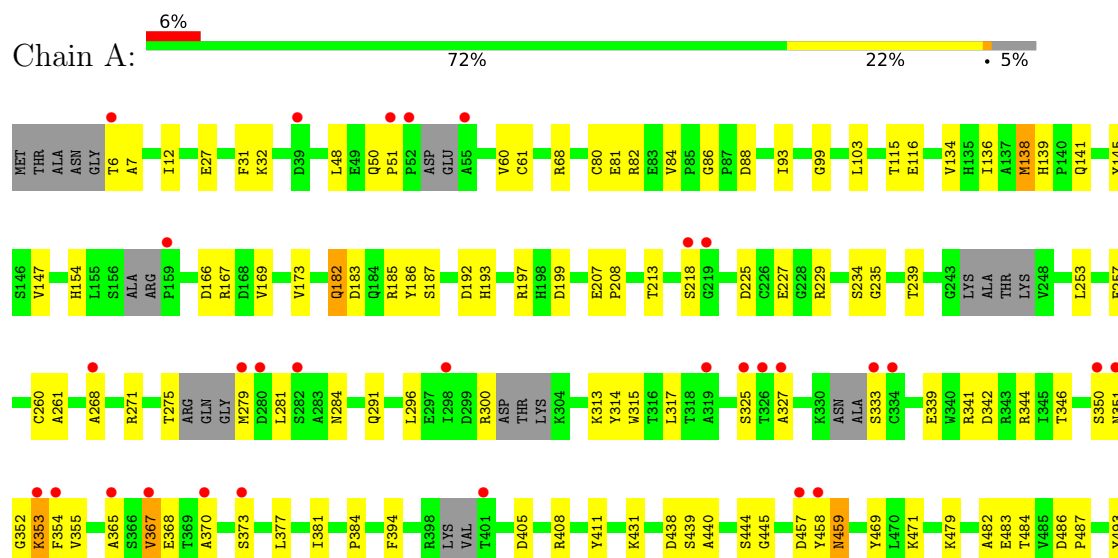
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	288	Total	O	0	0
			288	288		
6	B	273	Total	O	0	0
			273	273		

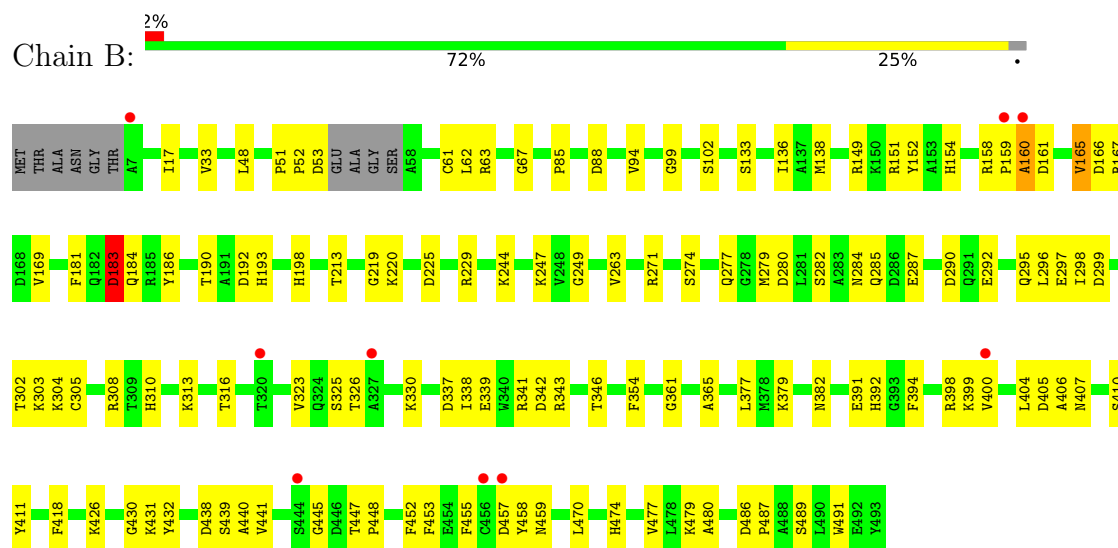
3 Residue-property plots

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: Fascin



• Molecule 1: Fascin



4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	161.25Å 70.72Å 111.06Å 90.00° 130.68° 90.00°	Depositor
Resolution (Å)	30.00 – 2.20 30.00 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.0 (30.00-2.20) 99.6 (30.00-2.20)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.81 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.196 , 0.234 0.200 , 0.235	Depositor DCC
R_{free} test set	2438 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å ²)	36.1	Xtriage
Anisotropy	0.197	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 58.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	0.026 for -h-2*k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8114	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 4.76% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CL, DTT, GOL, BR

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	0.57	0/3752	0.77	5/5066 (0.1%)
1	B	0.57	0/3883	0.79	3/5249 (0.1%)
All	All	0.57	0/7635	0.78	8/10315 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	160	ALA	N-CA-C	-9.64	97.89	110.33
1	A	367	VAL	N-CA-C	6.98	117.70	107.37
1	A	235	GLY	CA-C-N	6.79	126.58	119.05
1	A	235	GLY	C-N-CA	6.79	126.58	119.05
1	A	445	GLY	N-CA-C	5.66	118.96	111.03
1	A	116	GLU	CB-CA-C	-5.52	110.23	116.63
1	B	183	ASP	CB-CA-C	-5.48	109.77	117.23
1	B	382	ASN	N-CA-C	5.01	118.43	112.72

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	182	GLN	Peptide

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	3675	0	3540	94	0
1	B	3791	0	3676	98	0
2	A	2	0	0	2	0
2	B	4	0	0	1	0
3	A	5	0	0	1	0
3	B	8	0	0	0	0
4	A	18	0	24	2	0
4	B	42	0	56	10	0
5	A	8	0	10	2	0
6	A	288	0	0	6	0
6	B	273	0	0	2	0
All	All	8114	0	7306	191	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 13.

All (191) 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:B:160:ALA:HB1	1:B:161:ASP:CG	1.73	1.12
1:A:353:LYS:HE3	1:A:368:GLU:HG3	1.45	0.97
1:B:284:ASN:OD1	1:B:285:GLN:HG2	1.66	0.96
1:B:398:ARG:HG2	4:B:518:GOL:H11	1.52	0.89
1:A:260[B]:CYS:SG	1:A:458:TYR:OH	2.38	0.79
1:A:457:ASP:OD1	1:A:458:TYR:HB2	1.86	0.75
1:A:440:ALA:HA	1:A:479:LYS:HG2	1.71	0.73
1:A:218:SER:HB2	5:A:511:DTT:S1	2.31	0.70
1:A:260[B]:CYS:HG	1:A:458:TYR:HH	1.34	0.70
1:B:398:ARG:HG2	4:B:518:GOL:C1	2.23	0.69
1:A:145:TYR:HE2	4:A:510:GOL:H31	1.57	0.69
1:B:52:PRO:O	1:B:53:ASP:HB2	1.92	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:353:LYS:CE	1:A:368:GLU:HG3	2.22	0.69
1:B:159:PRO:C	1:B:160:ALA:O	2.34	0.67
1:B:313:LYS:HB3	1:B:325:SER:O	1.93	0.67
1:A:486:ASP:HB2	1:A:487:PRO:CD	2.24	0.66
1:A:253:LEU:HD11	5:A:511:DTT:H41	1.78	0.65
1:A:458:TYR:CG	1:A:459:ASN:N	2.57	0.65
1:B:99:GLY:HA2	1:B:213:THR:OG1	1.96	0.65
1:A:333:SER:HA	6:A:861:HOH:O	1.97	0.65
1:B:160:ALA:HB1	1:B:161:ASP:OD1	1.97	0.64
1:A:68:ARG:HD2	1:A:81:GLU:O	1.98	0.64
1:B:304:LYS:HE2	1:B:337:ASP:OD1	1.97	0.63
1:A:353:LYS:HD2	6:A:829:HOH:O	1.98	0.63
1:A:147:VAL:HA	4:A:510:GOL:H32	1.81	0.62
1:A:300:ARG:HG3	1:A:457:ASP:OD2	2.00	0.62
1:A:487:PRO:HD3	4:B:518:GOL:H12	1.82	0.62
1:A:86:GLY:HA3	2:A:502:BR:BR	2.54	0.61
1:A:183:ASP:N	1:A:183:ASP:OD1	2.32	0.61
1:B:399:LYS:N	4:B:518:GOL:H32	2.16	0.61
1:B:160:ALA:CB	1:B:161:ASP:CG	2.64	0.61
1:B:486:ASP:HB2	1:B:487:PRO:HD2	1.83	0.61
1:B:159:PRO:O	1:B:159:PRO:HG2	2.00	0.61
1:B:271:ARG:HD2	1:B:284:ASN:HA	1.83	0.60
1:A:48:LEU:HD11	1:A:60:VAL:HB	1.83	0.59
1:B:377:LEU:HD21	1:B:418:PHE:CZ	2.37	0.59
1:A:167:ARG:HD3	1:A:169:VAL:O	2.02	0.58
1:B:405:ASP:HB3	1:B:407:ASN:OD1	2.03	0.58
1:B:290:ASP:HB3	1:B:310:HIS:ND1	2.19	0.58
4:B:513:GOL:H31	6:B:863:HOH:O	2.03	0.58
1:A:197:ARG:NH2	1:A:199:ASP:OD2	2.37	0.58
1:B:160:ALA:HB1	1:B:161:ASP:CB	2.34	0.58
1:B:457:ASP:CG	1:B:458:TYR:H	2.12	0.57
1:B:486:ASP:H	1:B:489:SER:HG	1.53	0.57
1:A:377:LEU:C	1:A:377:LEU:HD23	2.30	0.56
1:B:160:ALA:CB	1:B:161:ASP:OD1	2.53	0.56
1:B:404:LEU:HD12	1:B:441:VAL:HG12	1.88	0.56
1:B:406:ALA:HB1	1:B:491:TRP:HH2	1.70	0.56
1:A:458:TYR:CD1	1:A:459:ASN:N	2.74	0.55
1:B:63:ARG:HH11	1:B:67:GLY:HA2	1.72	0.55
1:B:316:THR:O	1:B:323:VAL:HA	2.07	0.55
1:A:12:ILE:HB	1:A:48:LEU:HB3	1.89	0.54
1:B:277:GLN:O	1:B:280:ASP:HB2	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:391:GLU:HG3	1:B:392:HIS:CD2	2.42	0.54
1:A:182:GLN:O	1:A:185:ARG:N	2.41	0.54
1:B:339:GLU:OE1	1:B:341:ARG:NH1	2.41	0.54
1:A:484:THR:HG21	1:B:477:VAL:CG2	2.38	0.54
1:B:400:VAL:H	4:B:518:GOL:H32	1.73	0.54
1:A:341:ARG:O	1:A:342:ASP:HB2	2.07	0.53
1:B:354:PHE:O	1:B:365:ALA:HA	2.08	0.53
1:B:486:ASP:HB2	1:B:487:PRO:CD	2.38	0.53
1:A:31:PHE:HB3	1:A:80:CYS:O	2.08	0.53
1:A:353:LYS:HG3	1:A:367:VAL:O	2.08	0.53
1:A:82:ARG:HD3	3:A:504:CL:CL	2.46	0.52
1:A:350:SER:HB2	6:A:861:HOH:O	2.09	0.52
1:B:431:LYS:HD2	1:B:445:GLY:O	2.09	0.52
1:A:136:ILE:HG12	1:A:186:TYR:CZ	2.44	0.52
1:A:207:GLU:HB2	1:A:208:PRO:HD2	1.91	0.52
1:A:260[B]:CYS:SG	1:A:296:LEU:O	2.68	0.52
1:B:305[A]:CYS:SG	1:B:338:ILE:HD11	2.49	0.52
1:A:48:LEU:HD11	1:A:60:VAL:CB	2.40	0.51
1:B:17:ILE:HB	1:B:133:SER:HB2	1.91	0.51
1:B:94:VAL:HB	1:B:102[B]:SER:OG	2.10	0.51
1:B:406:ALA:HB1	1:B:491:TRP:CH2	2.44	0.51
1:B:279:MET:SD	1:B:326:THR:HG22	2.50	0.51
1:B:398:ARG:HE	1:B:405:ASP:CG	2.19	0.51
1:B:430:GLY:CA	4:B:515:GOL:H11	2.41	0.51
1:A:279:MET:HE1	1:A:313:LYS:NZ	2.26	0.50
1:A:279:MET:HE2	1:A:279:MET:HA	1.94	0.50
1:A:317:LEU:HD21	1:A:351:ASN:ND2	2.26	0.50
1:A:82:ARG:HD2	1:A:84:VAL:O	2.12	0.50
1:B:447:THR:HG23	1:B:448:PRO:HD2	1.92	0.50
1:A:6:THR:HG23	1:A:7:ALA:N	2.27	0.50
1:A:182:GLN:OE1	1:A:187:SER:HB3	2.12	0.50
1:B:343:ARG:HG2	1:B:452:PHE:HE1	1.77	0.49
1:B:470:LEU:CD2	1:B:480:ALA:HB2	2.43	0.49
1:B:426:LYS:HB2	1:B:431:LYS:O	2.12	0.49
1:A:394:PHE:CD1	1:A:411:TYR:HB3	2.48	0.49
1:A:115:THR:HB	6:A:617:HOH:O	2.12	0.49
1:A:457:ASP:HA	1:A:458:TYR:HB2	1.95	0.49
1:A:279:MET:HE1	1:A:313:LYS:HZ3	1.78	0.49
1:A:271:ARG:CZ	1:A:284:ASN:O	2.61	0.49
1:B:181:PHE:CE2	1:B:410:SER:HB3	2.48	0.48
1:B:159:PRO:O	1:B:159:PRO:CG	2.59	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:ASP:O	1:A:193:HIS:HB2	2.14	0.48
1:A:99:GLY:HA2	1:A:213:THR:OG1	2.14	0.48
1:B:52:PRO:HD3	1:B:85:PRO:HB2	1.96	0.48
1:B:440:ALA:HA	1:B:479:LYS:HG2	1.96	0.48
1:A:354:PHE:CE1	1:A:370:ALA:HB2	2.48	0.47
1:B:136:ILE:HG23	1:B:186:TYR:CZ	2.48	0.47
1:B:341:ARG:O	1:B:342:ASP:HB2	2.14	0.47
1:B:159:PRO:O	1:B:160:ALA:C	2.58	0.47
1:A:469:TYR:HB2	1:A:482:ALA:HB3	1.97	0.47
1:B:339:GLU:CD	1:B:341:ARG:HH11	2.22	0.47
1:A:139:HIS:CE1	1:A:141:GLN:HG3	2.49	0.47
1:A:93:ILE:HD13	1:A:103:LEU:HD23	1.97	0.47
1:A:268:ALA:CB	1:A:373:SER:HA	2.45	0.47
1:A:457:ASP:HA	1:A:458:TYR:CB	2.45	0.47
1:A:353:LYS:HG3	1:A:367:VAL:C	2.39	0.46
1:B:298:ILE:HD13	1:B:455:PHE:O	2.16	0.46
1:A:27:GLU:OE1	1:A:32:LYS:HB2	2.16	0.46
1:B:149:ARG:O	1:B:151:ARG:HG2	2.14	0.46
1:B:287:GLU:O	1:B:292:GLU:HG2	2.15	0.46
1:B:432:TYR:CD1	1:B:448:PRO:HB3	2.50	0.46
1:B:192:ASP:O	1:B:193:HIS:HB2	2.14	0.46
1:A:350:SER:CB	6:A:861:HOH:O	2.64	0.46
1:A:93:ILE:HD13	1:A:103:LEU:CD2	2.46	0.46
1:B:154:HIS:NE2	1:B:166:ASP:OD1	2.42	0.46
1:B:183:ASP:HB3	1:B:184:GLN:H	1.52	0.46
1:A:281:LEU:HD21	1:A:315:TRP:CE2	2.52	0.45
1:A:486:ASP:HB2	1:A:487:PRO:HD2	1.99	0.45
1:A:486:ASP:HB2	1:A:487:PRO:HD3	1.96	0.45
1:A:313:LYS:HA	1:A:327:ALA:O	2.17	0.45
1:B:296:LEU:HD12	1:B:305[B]:CYS:SG	2.56	0.45
1:B:198:HIS:CE1	1:B:229:ARG:HH22	2.35	0.45
1:A:355:VAL:HA	1:A:365:ALA:HA	1.99	0.45
1:B:377:LEU:HD21	1:B:418:PHE:HZ	1.81	0.45
1:B:457:ASP:CG	1:B:458:TYR:N	2.75	0.45
1:B:377:LEU:HD21	1:B:418:PHE:CE2	2.51	0.45
1:A:281:LEU:HD21	1:A:315:TRP:CD2	2.52	0.45
1:B:343:ARG:HD2	4:B:519:GOL:H32	1.99	0.45
1:A:225:ASP:OD1	1:A:225:ASP:C	2.60	0.44
1:A:341:ARG:HB2	1:A:344:ARG:O	2.17	0.44
1:B:302:THR:C	1:B:304:LYS:H	2.24	0.44
1:B:438:ASP:O	1:B:439:SER:HB2	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:ALA:HB3	1:A:257:GLU:OE1	2.17	0.44
1:B:88:ASP:HB2	2:B:501:BR:BR	2.71	0.44
1:B:61:CYS:C	1:B:62:LEU:HD12	2.43	0.44
1:A:136:ILE:HG23	1:A:138:MET:SD	2.57	0.44
1:B:296:LEU:CD1	1:B:305[B]:CYS:SG	3.06	0.44
1:A:234:SER:N	1:A:239:THR:O	2.49	0.44
1:B:470:LEU:HD23	1:B:480:ALA:HB2	1.99	0.44
1:A:483:GLU:OE1	1:B:474:HIS:CD2	2.71	0.44
1:B:190:THR:OG1	1:B:192:ASP:OD1	2.32	0.44
1:A:339:GLU:HB2	1:A:346:THR:OG1	2.18	0.43
1:A:351:ASN:OD1	1:A:352:GLY:N	2.51	0.43
1:B:263:VAL:HA	1:B:379:LYS:O	2.18	0.43
1:A:261:ALA:HB2	1:A:493:TYR:OH	2.18	0.43
1:B:158:ARG:HA	1:B:160:ALA:H	1.82	0.43
1:A:471:LYS:HD2	1:A:482:ALA:HB2	2.00	0.43
1:A:438:ASP:O	1:A:439:SER:HB2	2.18	0.43
1:B:274:SER:HB2	1:B:285:GLN:HG3	1.99	0.43
1:B:302:THR:O	1:B:303:LYS:HB2	2.17	0.43
1:A:6:THR:HG23	1:A:7:ALA:H	1.83	0.43
1:B:295:GLN:HB3	1:B:308:ARG:HB3	2.01	0.43
1:B:394:PHE:CD1	1:B:411:TYR:HB3	2.53	0.43
1:A:50:GLN:HG2	1:A:51:PRO:O	2.18	0.43
1:A:88:ASP:HB2	2:A:502:BR:BR	2.74	0.43
1:A:405:ASP:HB2	1:A:408:ARG:HD3	2.00	0.43
1:B:453:PHE:HB2	4:B:513:GOL:H32	2.01	0.43
1:B:158:ARG:HA	1:B:160:ALA:N	2.34	0.42
1:B:271:ARG:HH11	1:B:284:ASN:HA	1.84	0.42
1:B:299:ASP:O	1:B:303:LYS:N	2.49	0.42
1:A:173:VAL:HG23	1:A:384:PRO:HD3	2.02	0.42
1:B:459:ASN:OD1	1:B:459:ASN:N	2.49	0.42
1:B:219:GLY:C	1:B:220:LYS:HD2	2.44	0.42
1:B:152:TYR:C	1:B:165:VAL:HG13	2.45	0.42
1:B:160:ALA:HA	1:B:161:ASP:HA	1.50	0.42
1:B:48:LEU:HD12	1:B:48:LEU:HA	1.84	0.42
1:A:351:ASN:OD1	1:A:351:ASN:C	2.62	0.42
1:A:154:HIS:NE2	1:A:166:ASP:OD1	2.42	0.42
1:B:247:LYS:HB3	6:B:688:HOH:O	2.20	0.41
1:A:51:PRO:HB3	1:A:61:CYS:SG	2.61	0.41
1:A:353:LYS:HE3	1:A:368:GLU:HA	2.03	0.41
1:A:431:LYS:HD2	1:A:444:SER:O	2.21	0.41
1:B:225:ASP:OD2	1:B:229:ARG:NH2	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:354:PHE:HD2	1:A:367:VAL:O	2.04	0.41
1:A:381:ILE:C	1:A:381:ILE:HD12	2.46	0.41
1:B:51:PRO:O	1:B:52:PRO:C	2.64	0.41
1:B:244:LYS:NZ	1:B:249:GLY:O	2.54	0.41
1:A:192:ASP:O	1:A:193:HIS:CB	2.69	0.41
1:A:275:THR:HG22	1:A:291:GLN:O	2.20	0.41
1:A:314:TYR:O	1:A:325:SER:HA	2.21	0.41
1:B:342:ASP:O	4:B:513:GOL:H12	2.20	0.41
1:A:134:VAL:O	1:A:134:VAL:HG13	2.20	0.41
1:B:282:SER:HA	1:B:361:GLY:O	2.21	0.41
1:B:167:ARG:HD3	1:B:169:VAL:O	2.21	0.40
1:B:339:GLU:HB3	1:B:346:THR:OG1	2.21	0.40
1:A:50:GLN:NE2	6:A:821:HOH:O	2.54	0.40
1:A:227:GLU:HG3	1:A:229:ARG:NH2	2.37	0.40
1:B:297:GLU:OE1	1:B:330:LYS:HE3	2.21	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	455/493 (92%)	441 (97%)	14 (3%)	0	100	100
1	B	483/493 (98%)	467 (97%)	16 (3%)	0	100	100
All	All	938/986 (95%)	908 (97%)	30 (3%)	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	388/404 (96%)	385 (99%)	3 (1%)	73	85
1	B	401/404 (99%)	397 (99%)	4 (1%)	68	81
All	All	789/808 (98%)	782 (99%)	7 (1%)	70	84

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	138	MET
1	A	353	LYS
1	A	459	ASN
1	B	33	VAL
1	B	138	MET
1	B	165	VAL
1	B	183	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	GLN
1	A	77	ASN
1	A	266	GLN
1	A	392	HIS
1	B	277	GLN
1	B	392	HIS
1	B	415	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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 [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 30 ligands modelled in this entry, 19 are monoatomic - leaving 11 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	$\# Z > 2$
4	GOL	A	510	-	5,5,5	0.23	0	5,5,5	0.36	0
4	GOL	B	514	-	5,5,5	0.27	0	5,5,5	0.24	0
4	GOL	B	519	-	5,5,5	0.24	0	5,5,5	0.42	0
4	GOL	A	509	-	5,5,5	0.30	0	5,5,5	0.15	0
4	GOL	B	516	-	5,5,5	0.29	0	5,5,5	0.38	0
4	GOL	B	518	-	5,5,5	0.37	0	5,5,5	0.65	0
4	GOL	B	515	-	5,5,5	0.26	0	5,5,5	0.38	0
4	GOL	B	513	-	5,5,5	0.25	0	5,5,5	0.57	0
4	GOL	A	508	-	5,5,5	0.11	0	5,5,5	0.58	0
4	GOL	B	517	-	5,5,5	0.28	0	5,5,5	0.43	0
5	DTT	A	511	-	7,7,7	0.36	0	4,8,8	1.95	1 (25%)

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
4	GOL	A	510	-	-	4/4/4/4	-
4	GOL	B	514	-	-	0/4/4/4	-
4	GOL	B	519	-	-	2/4/4/4	-
4	GOL	A	509	-	-	1/4/4/4	-
4	GOL	B	516	-	-	2/4/4/4	-
4	GOL	B	518	-	-	2/4/4/4	-
4	GOL	B	515	-	-	4/4/4/4	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	GOL	B	513	-	-	1/4/4/4	-
4	GOL	A	508	-	-	2/4/4/4	-
4	GOL	B	517	-	-	2/4/4/4	-
5	DTT	A	511	-	-	3/8/8/8	-

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
5	A	511	DTT	C3-C4-S4	-3.69	104.12	114.43

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	508	GOL	O1-C1-C2-C3
4	B	515	GOL	O1-C1-C2-C3
4	B	516	GOL	C1-C2-C3-O3
4	B	516	GOL	O2-C2-C3-O3
4	B	517	GOL	O1-C1-C2-C3
4	B	519	GOL	O1-C1-C2-C3
5	A	511	DTT	C2-C3-C4-S4
5	A	511	DTT	O3-C3-C4-S4
4	A	508	GOL	O1-C1-C2-O2
4	B	517	GOL	O1-C1-C2-O2
4	A	509	GOL	O1-C1-C2-C3
4	A	510	GOL	C1-C2-C3-O3
4	B	518	GOL	C1-C2-C3-O3
4	B	519	GOL	O1-C1-C2-O2
4	B	515	GOL	O1-C1-C2-O2
4	B	515	GOL	C1-C2-C3-O3
4	A	510	GOL	O1-C1-C2-O2
5	A	511	DTT	S1-C1-C2-O2
4	A	510	GOL	O1-C1-C2-C3
4	B	518	GOL	O2-C2-C3-O3
4	A	510	GOL	O2-C2-C3-O3
4	B	515	GOL	O2-C2-C3-O3
4	B	513	GOL	C1-C2-C3-O3

There are no ring outliers.

6 monomers are involved in 14 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	510	GOL	2	0
4	B	519	GOL	1	0
4	B	518	GOL	5	0
4	B	515	GOL	1	0
4	B	513	GOL	3	0
5	A	511	DTT	2	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 ⓘ

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	470/493 (95%)	0.27	30 (6%) 25 22	28, 48, 78, 98	2 (0%)
1	B	483/493 (97%)	0.14	9 (1%) 66 63	24, 47, 74, 88	4 (0%)
All	All	953/986 (96%)	0.20	39 (4%) 41 38	24, 48, 76, 98	6 (0%)

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	160	ALA	4.3
1	B	444	SER	3.5
1	A	370	ALA	3.5
1	A	457	ASP	3.2
1	B	327	ALA	3.1
1	A	334	CYS	3.1
1	A	52	PRO	3.1
1	A	268	ALA	3.0
1	B	159	PRO	3.0
1	A	219	GLY	2.8
1	A	365	ALA	2.8
1	A	218	SER	2.7
1	A	373	SER	2.7
1	A	367	VAL	2.7
1	B	400	VAL	2.7
1	A	51	PRO	2.7
1	A	458	TYR	2.7
1	A	55	ALA	2.7
1	A	351	ASN	2.7
1	B	320	THR	2.6
1	A	333	SER	2.6
1	A	280	ASP	2.6
1	A	319	ALA	2.5
1	A	327	ALA	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	326	THR	2.4
1	A	298	ILE	2.4
1	B	457	ASP	2.4
1	A	350	SER	2.4
1	A	325	SER	2.3
1	A	39	ASP	2.3
1	B	456	CYS	2.3
1	A	159	PRO	2.2
1	B	7	ALA	2.2
1	A	354	PHE	2.2
1	A	401	THR	2.2
1	A	279	MET	2.1
1	A	353	LYS	2.1
1	A	282	SER	2.0
1	A	6	THR	2.0

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
5	DTT	A	511	8/8	0.80	0.18	79,84,85,87	0
4	GOL	B	518	6/6	0.81	0.14	44,47,48,51	0
4	GOL	B	519	6/6	0.83	0.13	77,78,79,79	0
4	GOL	B	515	6/6	0.83	0.13	64,65,66,69	0
4	GOL	B	516	6/6	0.86	0.14	66,67,68,68	0
4	GOL	A	510	6/6	0.87	0.11	56,57,59,61	0
4	GOL	B	517	6/6	0.88	0.11	59,62,64,66	0
3	CL	B	505	1/1	0.89	0.12	57,57,57,57	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CL	A	504	1/1	0.90	0.17	67,67,67,67	0
3	CL	B	512	1/1	0.92	0.13	70,70,70,70	0
3	CL	A	505	1/1	0.92	0.12	66,66,66,66	0
2	BR	B	504	1/1	0.92	0.16	86,86,86,86	0
3	CL	B	508	1/1	0.92	0.10	62,62,62,62	0
4	GOL	A	508	6/6	0.93	0.10	42,47,51,53	0
3	CL	A	503	1/1	0.93	0.08	57,57,57,57	0
4	GOL	B	513	6/6	0.93	0.13	43,46,48,49	0
3	CL	B	509	1/1	0.94	0.15	56,56,56,56	0
3	CL	A	507	1/1	0.95	0.07	54,54,54,54	0
3	CL	B	506	1/1	0.95	0.08	65,65,65,65	0
2	BR	A	502	1/1	0.96	0.07	57,57,57,57	0
4	GOL	A	509	6/6	0.96	0.08	41,45,46,49	0
4	GOL	B	514	6/6	0.96	0.07	39,41,42,43	0
2	BR	B	503	1/1	0.97	0.06	75,75,75,75	0
3	CL	A	506	1/1	0.98	0.11	36,36,36,36	0
2	BR	B	502	1/1	0.98	0.05	75,75,75,75	0
2	BR	A	501	1/1	0.98	0.09	75,75,75,75	0
3	CL	B	511	1/1	0.98	0.06	44,44,44,44	0
3	CL	B	507	1/1	0.99	0.09	37,37,37,37	0
3	CL	B	510	1/1	0.99	0.09	34,34,34,34	0
2	BR	B	501	1/1	0.99	0.04	50,50,50,50	0

6.5 Other polymers [i](#)

There are no such residues in this entry.