



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 08:06 PM UTC

PDB ID : 4GOY / pdb_00004goy
Title : The crystal structure of human fascin 1 K41A mutant
Authors : Yang, S.Y.; Huang, F.K.; Huang, J.; Chen, S.; Jakoncic, J.; Leo-Macias, A.;
Diaz-Avalos, R.; Chen, L.; Zhang, J.J.; Huang, X.Y.
Deposited on : 2012-08-20
Resolution : 2.30 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

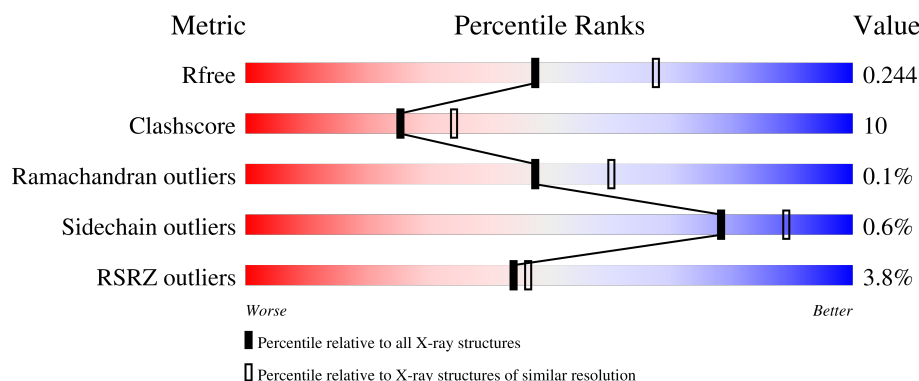
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

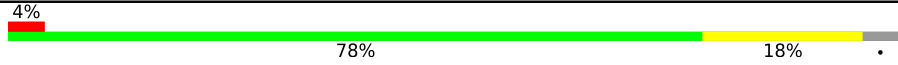
The reported resolution of this entry is 2.30 Å.

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	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	
1	B	493	

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	501	-	-	X	-
2	BR	B	504	-	-	X	-

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 8000 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	475	Total	C	N	O	S	0	1	0
			3701	2317	656	713	15			
1	B	484	Total	C	N	O	S	0	3	0
			3786	2367	676	726	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	41	ALA	LYS	engineered mutation	UNP Q16658
B	41	ALA	LYS	engineered mutation	UNP Q16658

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

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	3	Total	Br	0	0
			3	3		
2	B	5	Total	Br	0	0
			5	5		

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

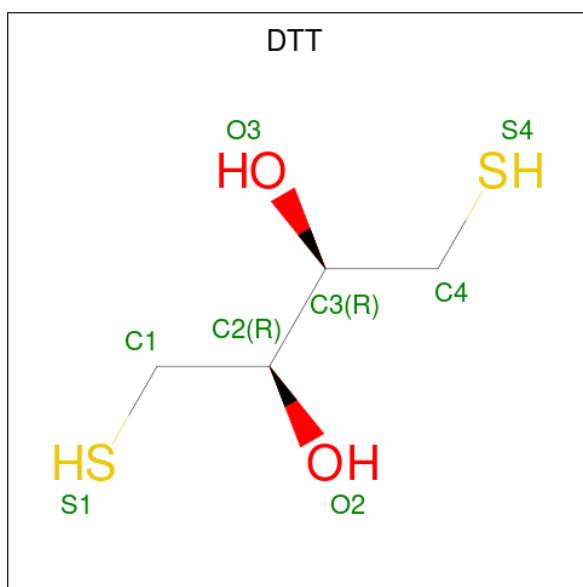
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	4	Total	Cl	0	0
			4	4		
3	B	6	Total	Cl	0	0
			6	6		

- 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		

- 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	B	1	Total	C	O	S	0	0
			8	4	2	2		

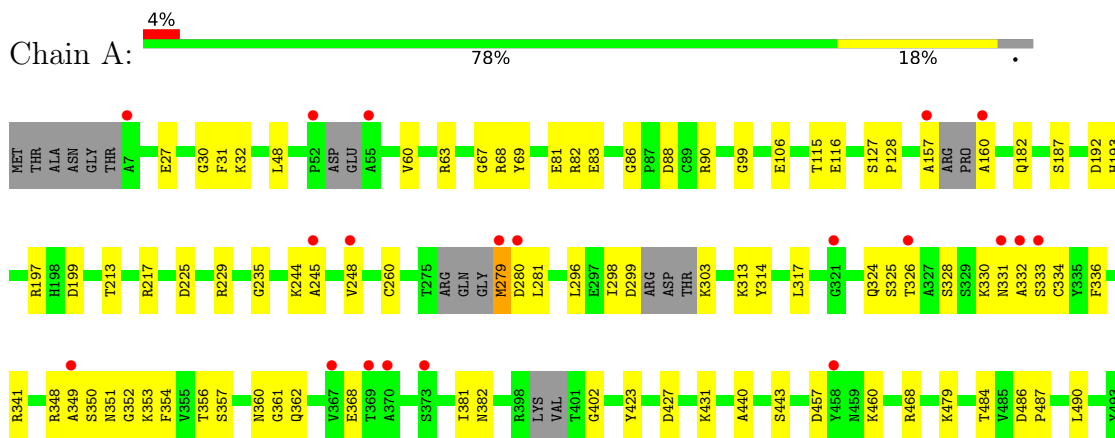
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	222	Total	O	0	0
			222	222		
6	B	229	Total	O	0	0
			229	229		

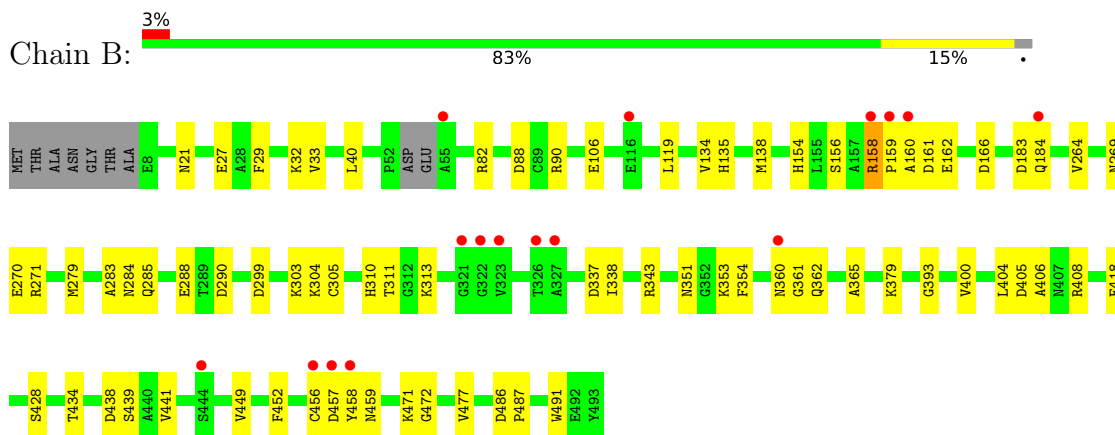
3 Residue-property plots

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	161.14Å 70.59Å 110.98Å 90.00° 130.63° 90.00°	Depositor
Resolution (Å)	30.00 – 2.30 30.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	98.8 (30.00-2.30) 99.7 (30.00-2.30)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.17 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.201 , 0.245 0.201 , 0.244	Depositor DCC
R_{free} test set	2129 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	35.8	Xtriage
Anisotropy	0.194	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 53.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.025 for -h-2*1,-k,l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	8000	wwPDB-VP
Average B, all atoms (Å ²)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.81% 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: DTT, BR, GOL, CL

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	0.71	0/3779	0.84	4/5105 (0.1%)
1	B	0.70	0/3875	0.88	4/5239 (0.1%)
All	All	0.71	0/7654	0.86	8/10344 (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
1	B	0	1
All	All	0	2

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	162	GLU	N-CA-C	-6.28	99.78	109.14
1	B	361	GLY	N-CA-C	-5.48	107.10	115.00
1	A	235	GLY	CA-C-N	5.35	125.16	119.28
1	A	235	GLY	C-N-CA	5.35	125.16	119.28
1	B	160	ALA	CA-C-N	-5.34	113.53	122.17
1	B	160	ALA	C-N-CA	-5.34	113.53	122.17
1	A	116	GLU	CB-CA-C	-5.27	110.48	116.54
1	A	402	GLY	N-CA-C	-5.06	108.63	115.21

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	244	LYS	Peptide
1	B	456	CYS	Peptide

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	3701	0	3574	86	0
1	B	3786	0	3673	57	0
2	A	3	0	0	2	0
2	B	5	0	0	3	0
3	A	4	0	0	0	0
3	B	6	0	0	0	0
4	A	18	0	24	3	0
4	B	18	0	24	2	0
5	B	8	0	10	1	0
6	A	222	0	0	12	0
6	B	229	0	0	4	0
All	All	8000	0	7305	142	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 10.

All (142) 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:360:ASN:HD22	1:B:362:GLN:NE2	1.51	1.09
1:B:360:ASN:ND2	1:B:362:GLN:HE21	1.56	1.03
1:A:48:LEU:HD11	1:A:60:VAL:HB	1.41	0.99
1:B:29:PHE:HB2	5:B:515:DTT:H42	1.52	0.90
1:A:68:ARG:NH1	1:A:81:GLU:O	2.06	0.87
1:A:348:ARG:CD	6:A:783:HOH:O	2.22	0.86
1:A:299:ASP:HB2	1:A:303:LYS:N	1.92	0.85
1:B:279:MET:HE2	1:B:313:LYS:HZ3	1.43	0.83
1:A:341:ARG:HG2	1:A:341:ARG:HH11	1.47	0.79
1:A:30:GLY:O	1:A:32:LYS:HG3	1.81	0.79
1:A:48:LEU:HD11	1:A:60:VAL:CB	2.12	0.79

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:30:GLY:C	1:A:32:LYS:HG3	2.10	0.76
1:A:182:GLN:NE2	1:A:187:SER:OG	2.21	0.73
1:A:324:GLN:HG3	1:A:326:THR:OG1	1.88	0.73
1:B:82:ARG:HD3	2:B:504:BR:BR	2.44	0.72
1:A:348:ARG:CG	6:A:783:HOH:O	2.38	0.72
1:B:290:ASP:HB3	1:B:310:HIS:ND1	2.06	0.71
1:A:353:LYS:HE2	1:A:368:GLU:HG2	1.73	0.70
1:A:486:ASP:HB2	1:A:487:PRO:CD	2.21	0.70
1:B:279:MET:HE2	1:B:313:LYS:NZ	2.06	0.70
1:A:348:ARG:HD3	6:A:783:HOH:O	1.86	0.70
1:A:356:THR:HG22	1:A:357:SER:N	2.07	0.70
1:A:349:ALA:O	1:A:352:GLY:N	2.13	0.67
1:B:351:ASN:HB3	1:B:353:LYS:H	1.59	0.67
1:A:299:ASP:CB	1:A:303:LYS:N	2.56	0.67
1:A:68:ARG:CZ	1:A:83:GLU:HG3	2.25	0.67
1:B:459:ASN:OD1	1:B:459:ASN:N	2.29	0.66
1:A:353:LYS:HE2	1:A:368:GLU:CG	2.26	0.66
1:B:393:GLY:HA2	4:B:513:GOL:H12	1.79	0.65
1:A:486:ASP:HB2	1:A:487:PRO:HD3	1.78	0.65
1:A:48:LEU:CD1	1:A:60:VAL:HB	2.23	0.64
1:A:115:THR:HB	6:A:668:HOH:O	1.97	0.64
1:B:284:ASN:OD1	1:B:285:GLN:HG2	1.98	0.63
1:B:158:ARG:H	1:B:159:PRO:CD	2.11	0.63
1:A:360:ASN:C	1:A:360:ASN:OD1	2.42	0.62
1:A:82:ARG:HG2	6:A:767:HOH:O	1.98	0.62
1:B:158:ARG:H	1:B:159:PRO:HD2	1.63	0.62
1:A:30:GLY:CA	1:A:32:LYS:HE2	2.30	0.62
1:A:423:TYR:HE2	4:A:508:GOL:H31	1.65	0.62
1:A:431:LYS:HE2	1:A:443:SER:O	2.00	0.61
1:A:314:TYR:CZ	1:A:330:LYS:HE2	2.35	0.61
1:A:341:ARG:HG2	1:A:341:ARG:NH1	2.11	0.61
1:A:99:GLY:HA2	1:A:213:THR:OG1	2.00	0.61
1:B:360:ASN:HB2	1:B:362:GLN:HG2	1.82	0.60
1:A:331:ASN:HB2	6:A:624:HOH:O	2.01	0.60
1:A:317:LEU:HB3	1:A:333:SER:HB3	1.83	0.60
1:A:348:ARG:HG3	6:A:783:HOH:O	1.99	0.60
1:A:127:SER:HB2	1:A:128:PRO:HD2	1.84	0.59
1:A:354:PHE:CG	6:A:783:HOH:O	2.52	0.59
1:B:82:ARG:CD	2:B:504:BR:BR	3.05	0.59
1:A:63:ARG:HE	1:A:67:GLY:HA2	1.67	0.59
1:B:351:ASN:ND2	1:B:353:LYS:HD3	2.18	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:458:TYR:HB3	1:B:459:ASN:OD1	2.04	0.58
1:B:158:ARG:HB2	1:B:159:PRO:HD3	1.86	0.57
1:A:440:ALA:HA	1:A:479:LYS:HG2	1.86	0.57
1:B:310:HIS:CD2	1:B:311:THR:HG23	2.39	0.57
1:B:487:PRO:HD3	6:B:803:HOH:O	2.02	0.57
1:B:134:VAL:HG11	6:B:766:HOH:O	2.05	0.57
1:A:332:ALA:O	1:A:350:SER:HB2	2.05	0.56
1:A:27:GLU:OE1	1:A:32:LYS:HB2	2.06	0.56
1:A:217:ARG:NH1	1:A:248:VAL:HB	2.20	0.56
1:B:406:ALA:HB1	1:B:491:TRP:HH2	1.71	0.56
1:A:88:ASP:HB2	2:A:501:BR:BR	2.60	0.56
1:A:484:THR:HG21	1:B:477:VAL:CG2	2.37	0.55
1:A:30:GLY:HA3	1:A:32:LYS:HE2	1.87	0.55
1:B:264:VAL:HG21	1:B:288:GLU:HG2	1.89	0.54
1:B:158:ARG:N	1:B:159:PRO:CD	2.70	0.54
1:A:351:ASN:OD1	1:A:353:LYS:HB2	2.08	0.54
1:A:331:ASN:H	1:A:334:CYS:HB2	1.72	0.54
1:A:229:ARG:HA	1:A:245:ALA:O	2.08	0.54
1:A:353:LYS:CE	1:A:368:GLU:HG2	2.37	0.54
1:A:317:LEU:CD2	1:A:351:ASN:HB2	2.38	0.53
1:A:281:LEU:HD21	1:A:325:SER:HB3	1.90	0.52
1:A:328:SER:H	4:A:510:GOL:H32	1.74	0.52
1:B:269:ASN:HD21	1:B:283:ALA:HB3	1.74	0.52
1:A:68:ARG:CZ	1:A:83:GLU:CG	2.87	0.52
1:A:457:ASP:HB3	1:A:460:LYS:HB2	1.92	0.52
1:B:183:ASP:O	1:B:184:GLN:CB	2.56	0.52
1:A:86:GLY:HA3	2:A:501:BR:BR	2.64	0.52
1:A:260[B]:CYS:SG	1:A:296:LEU:O	2.68	0.51
1:A:317:LEU:HD21	1:A:351:ASN:HB2	1.92	0.50
1:B:27:GLU:HB2	1:B:32:LYS:HB2	1.94	0.50
1:A:490:LEU:HD13	1:B:400:VAL:HG11	1.94	0.50
1:A:354:PHE:N	6:A:783:HOH:O	2.45	0.49
1:B:360:ASN:ND2	1:B:362:GLN:NE2	2.32	0.49
1:A:68:ARG:NH1	1:A:83:GLU:HG3	2.27	0.49
1:A:298:ILE:O	1:A:299:ASP:C	2.54	0.49
1:B:354:PHE:O	1:B:365:ALA:HA	2.12	0.49
1:B:88:ASP:HB2	2:B:501:BR:BR	2.69	0.48
1:B:304:LYS:NZ	1:B:337:ASP:OD1	2.38	0.48
1:B:360:ASN:HD22	1:B:362:GLN:HE21	0.71	0.48
1:A:361:GLY:O	1:A:362:GLN:C	2.55	0.48
1:B:271:ARG:HD3	1:B:284:ASN:HA	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:63:ARG:HD2	1:A:69:TYR:CZ	2.49	0.47
1:A:63:ARG:NE	1:A:67:GLY:HA2	2.30	0.47
1:A:354:PHE:CD1	6:A:783:HOH:O	2.67	0.47
1:A:356:THR:CG2	1:A:357:SER:N	2.74	0.47
1:B:90:ARG:HD3	1:B:106:GLU:OE1	2.15	0.47
1:B:404:LEU:HD12	1:B:441:VAL:HG12	1.97	0.47
1:A:330:LYS:HE3	6:A:762:HOH:O	2.15	0.47
1:A:427:ASP:OD1	1:A:431:LYS:HG2	2.15	0.47
1:A:30:GLY:HA3	1:A:32:LYS:CE	2.46	0.46
1:A:336:PHE:CE1	1:A:349:ALA:HB2	2.50	0.46
1:B:183:ASP:O	1:B:184:GLN:HB3	2.16	0.46
1:B:279:MET:CE	1:B:313:LYS:HZ3	2.23	0.46
1:B:299:ASP:O	1:B:303:LYS:N	2.46	0.45
1:A:423:TYR:HE2	4:A:508:GOL:C3	2.28	0.45
1:B:310:HIS:HD2	1:B:311:THR:HG23	1.81	0.45
1:A:333:SER:HB2	6:A:624:HOH:O	2.17	0.44
1:B:303:LYS:HG2	6:B:772:HOH:O	2.17	0.44
1:B:486:ASP:HB2	1:B:487:PRO:HD2	2.00	0.44
1:A:90:ARG:HD3	1:A:106:GLU:OE1	2.17	0.44
1:A:68:ARG:HD2	1:A:81:GLU:O	2.17	0.44
1:B:343:ARG:HG2	1:B:452:PHE:HE1	1.83	0.44
1:B:379:LYS:HD3	1:B:418:PHE:CD2	2.52	0.44
1:A:31:PHE:CD1	1:A:81:GLU:HG3	2.53	0.43
1:B:154:HIS:NE2	1:B:166:ASP:OD1	2.48	0.43
1:A:279:MET:HE1	1:A:313:LYS:HG3	2.00	0.43
1:A:353:LYS:CD	1:A:368:GLU:HG2	2.49	0.43
1:A:280:ASP:OD2	1:A:280:ASP:N	2.48	0.43
1:A:331:ASN:O	1:A:334:CYS:HB2	2.19	0.43
1:A:486:ASP:CB	1:A:487:PRO:CD	2.95	0.43
1:B:269:ASN:O	1:B:270:GLU:HB2	2.19	0.42
1:A:381:ILE:HD12	1:A:382:ASN:N	2.34	0.42
1:B:305[B]:CYS:SG	1:B:338:ILE:HD11	2.58	0.42
1:B:471:LYS:HD3	1:B:472:GLY:O	2.18	0.42
1:B:434:THR:HG22	1:B:449:VAL:HG23	2.02	0.42
1:B:438:ASP:O	1:B:439:SER:HB2	2.19	0.42
1:A:225:ASP:OD1	1:A:225:ASP:C	2.62	0.41
1:B:405:ASP:HB2	1:B:408:ARG:HD3	2.02	0.41
1:B:158:ARG:N	1:B:159:PRO:HD2	2.31	0.41
1:A:30:GLY:HA3	1:A:32:LYS:CD	2.51	0.41
1:A:192:ASP:O	1:A:193:HIS:HB2	2.21	0.41
1:A:325:SER:O	1:A:325:SER:OG	2.37	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:360:ASN:OD1	1:A:360:ASN:O	2.38	0.41
1:B:457:ASP:OD1	1:B:458:TYR:N	2.54	0.41
1:B:33:VAL:HG12	1:B:119:LEU:HD12	2.03	0.41
1:B:40:LEU:HD22	1:B:135:HIS:CG	2.56	0.41
1:A:197:ARG:NH1	1:A:199:ASP:OD2	2.53	0.40
4:B:513:GOL:H11	6:B:679:HOH:O	2.19	0.40
1:B:156:SER:O	1:B:161:ASP:OD1	2.38	0.40
1:A:157:ALA:C	1:A:160:ALA:O	2.64	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	464/493 (94%)	440 (95%)	24 (5%)	0	100	100
1	B	483/493 (98%)	458 (95%)	24 (5%)	1 (0%)	43	55
All	All	947/986 (96%)	898 (95%)	48 (5%)	1 (0%)	48	60

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	158	ARG

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	389/403 (96%)	387 (100%)	2 (0%)	81	90
1	B	400/403 (99%)	397 (99%)	3 (1%)	73	86
All	All	789/806 (98%)	784 (99%)	5 (1%)	78	89

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	279	MET
1	A	468	ARG
1	B	21	ASN
1	B	138	MET
1	B	428	SER

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	65	HIS
1	A	143	ASN
1	A	182	GLN
1	A	198	HIS
1	A	459	ASN
1	B	331	ASN
1	B	362	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 25 ligands modelled in this entry, 18 are monoatomic - leaving 7 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	B	513	-	5,5,5	0.32	0	5,5,5	0.34	0
4	GOL	A	509	-	5,5,5	0.26	0	5,5,5	0.81	0
4	GOL	B	514	-	5,5,5	0.21	0	5,5,5	0.46	0
4	GOL	B	512	-	5,5,5	0.28	0	5,5,5	0.53	0
4	GOL	A	508	-	5,5,5	0.23	0	5,5,5	0.49	0
5	DTT	B	515	-	7,7,7	0.54	0	4,8,8	1.27	0
4	GOL	A	510	-	5,5,5	0.23	0	5,5,5	0.56	0

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	B	513	-	-	2/4/4/4	-
4	GOL	A	509	-	-	3/4/4/4	-
4	GOL	B	514	-	-	2/4/4/4	-
4	GOL	B	512	-	-	0/4/4/4	-
4	GOL	A	508	-	-	4/4/4/4	-
5	DTT	B	515	-	-	4/8/8/8	-
4	GOL	A	510	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (17) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	508	GOL	O1-C1-C2-O2
4	A	508	GOL	O1-C1-C2-C3
4	A	509	GOL	O1-C1-C2-C3
4	A	510	GOL	O2-C2-C3-O3

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Mol	Chain	Res	Type	Atoms
5	B	515	DTT	C1-C2-C3-O3
5	B	515	DTT	C1-C2-C3-C4
5	B	515	DTT	O2-C2-C3-O3
5	B	515	DTT	O2-C2-C3-C4
4	A	508	GOL	C1-C2-C3-O3
4	A	510	GOL	C1-C2-C3-O3
4	B	513	GOL	C1-C2-C3-O3
4	B	514	GOL	C1-C2-C3-O3
4	A	508	GOL	O2-C2-C3-O3
4	A	509	GOL	O1-C1-C2-O2
4	B	514	GOL	O2-C2-C3-O3
4	B	513	GOL	O2-C2-C3-O3
4	A	509	GOL	C1-C2-C3-O3

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	513	GOL	2	0
4	A	508	GOL	2	0
5	B	515	DTT	1	0
4	A	510	GOL	1	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	475/493 (96%)	0.09	20 (4%) 40 42	19, 36, 67, 89	1 (0%)
1	B	484/493 (98%)	0.02	16 (3%) 49 51	16, 34, 65, 90	3 (0%)
All	All	959/986 (97%)	0.05	36 (3%) 44 46	16, 35, 67, 90	4 (0%)

All (36) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	457	ASP	5.6
1	B	160	ALA	4.8
1	A	160	ALA	3.9
1	A	332	ALA	3.8
1	A	370	ALA	3.7
1	B	458	TYR	3.4
1	A	458	TYR	3.3
1	A	157	ALA	3.0
1	B	158	ARG	2.9
1	A	349	ALA	2.9
1	A	369	THR	2.9
1	B	321	GLY	2.7
1	B	456	CYS	2.7
1	B	327	ALA	2.7
1	B	55	ALA	2.7
1	B	159	PRO	2.7
1	B	444	SER	2.6
1	A	280	ASP	2.5
1	A	55	ALA	2.4
1	A	245	ALA	2.4
1	B	326	THR	2.4
1	A	333	SER	2.3
1	A	326	THR	2.3
1	A	367	VAL	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	116	GLU	2.3
1	A	321	GLY	2.2
1	A	373	SER	2.2
1	B	360	ASN	2.1
1	A	248	VAL	2.1
1	B	322	GLY	2.1
1	B	184	GLN	2.1
1	A	52	PRO	2.1
1	B	323	VAL	2.1
1	A	279	MET	2.0
1	A	7	ALA	2.0
1	A	331	ASN	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
3	CL	A	507	1/1	0.76	0.19	62,62,62,62	0
4	GOL	A	510	6/6	0.83	0.16	58,62,63,63	0
5	DTT	B	515	8/8	0.83	0.18	61,66,69,77	0
3	CL	B	509	1/1	0.85	0.16	74,74,74,74	0
4	GOL	A	509	6/6	0.87	0.12	33,40,43,47	0
4	GOL	B	512	6/6	0.90	0.10	28,32,35,38	0
3	CL	B	511	1/1	0.90	0.12	48,48,48,48	0
3	CL	A	504	1/1	0.91	0.09	48,48,48,48	0
4	GOL	A	508	6/6	0.91	0.12	30,38,40,42	0
4	GOL	B	514	6/6	0.92	0.12	35,39,41,44	0
3	CL	A	506	1/1	0.93	0.10	58,58,58,58	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	GOL	B	513	6/6	0.93	0.12	39,42,46,47	0
3	CL	B	508	1/1	0.95	0.08	44,44,44,44	0
2	BR	A	501	1/1	0.96	0.05	49,49,49,49	0
3	CL	B	507	1/1	0.96	0.11	43,43,43,43	0
2	BR	B	504	1/1	0.96	0.08	65,65,65,65	0
2	BR	A	503	1/1	0.97	0.08	65,65,65,65	0
2	BR	A	502	1/1	0.97	0.07	49,49,49,49	0
2	BR	B	505	1/1	0.97	0.05	66,66,66,66	0
3	CL	B	510	1/1	0.97	0.08	63,63,63,63	0
3	CL	A	505	1/1	0.98	0.06	41,41,41,41	0
2	BR	B	503	1/1	0.98	0.09	52,52,52,52	0
2	BR	B	501	1/1	0.98	0.04	41,41,41,41	0
3	CL	B	506	1/1	0.98	0.05	35,35,35,35	0
2	BR	B	502	1/1	0.99	0.06	45,45,45,45	0

6.5 Other polymers [i](#)

There are no such residues in this entry.