



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 1, 2026 – 04:08 PM UTC

PDB ID : 4GP3 / pdb_00004gp3
Title : The crystal structure of human fascin 1 K358A 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 : 2.25 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

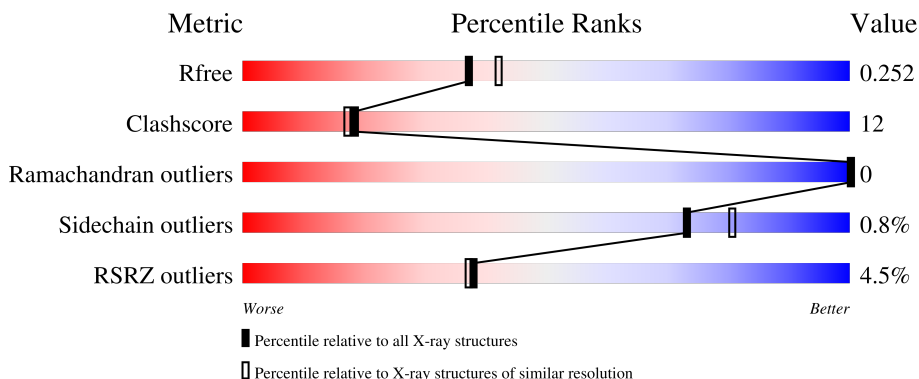
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	4% 75% 22% .
1	B	493	4% 81% 16% .

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	-
3	CL	B	507	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7976 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	480	Total	C	N	O	S	0	2	0
			3746	2344	665	721	16			
1	B	481	Total	C	N	O	S	0	3	0
			3764	2354	671	722	17			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	358	ALA	LYS	engineered mutation	UNP Q16658
B	358	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	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	13	Total	Cl	0	0
			13	13		

- 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		

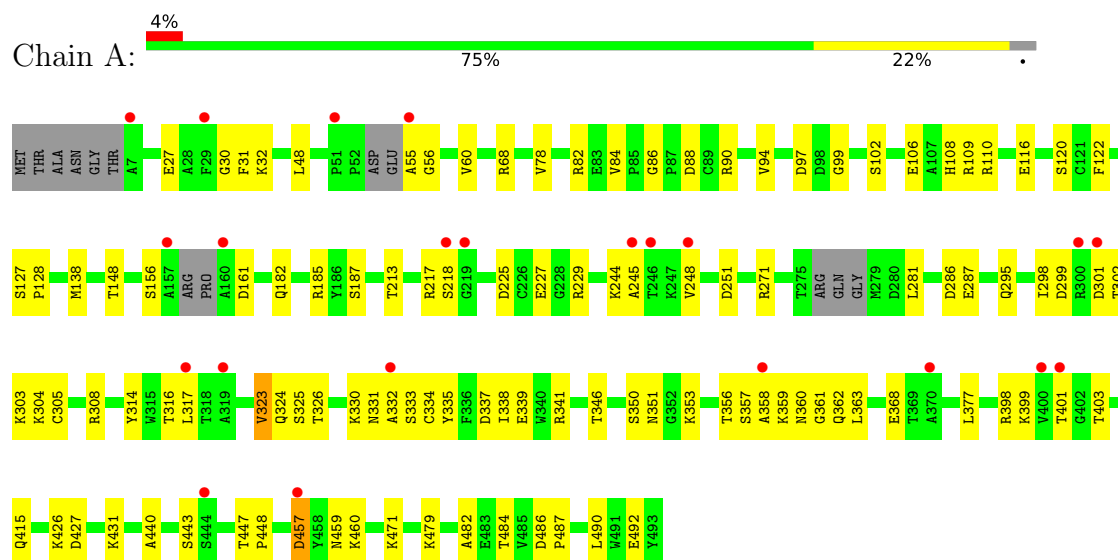
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	199	Total	O	0	0
			199	199		
5	B	200	Total	O	0	0
			200	200		

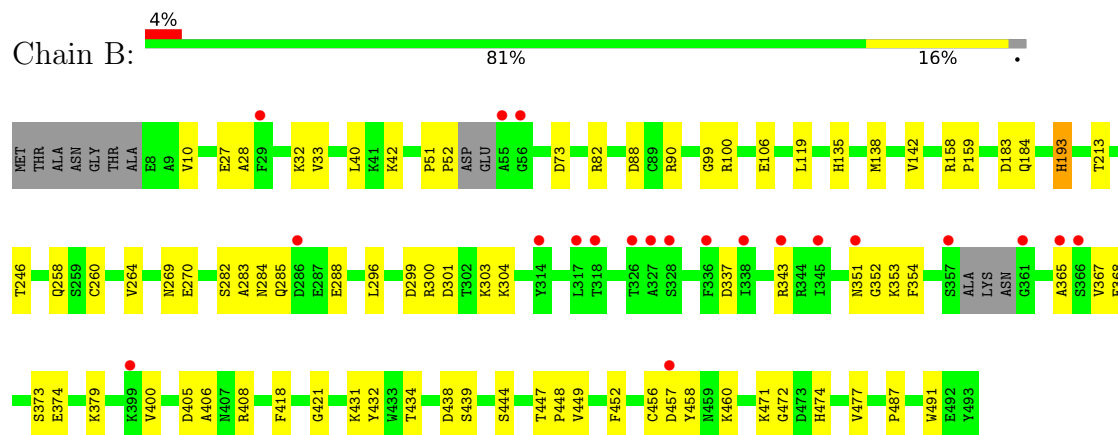
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: 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	160.91Å 70.74Å 110.51Å 90.00° 130.65° 90.00°	Depositor
Resolution (Å)	30.00 – 2.25 30.00 – 2.25	Depositor EDS
% Data completeness (in resolution range)	99.2 (30.00-2.25) 99.2 (30.00-2.25)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.68 (at 2.24Å)	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
R, R_{free}	0.203 , 0.252 0.203 , 0.252	Depositor DCC
R_{free} test set	2256 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	38.9	Xtriage
Anisotropy	0.562	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 47.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.022 for -h-2*1,-k,l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7976	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.82% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.65	0/3829	0.79	3/5175 (0.1%)
1	B	0.64	0/3852	0.82	1/5207 (0.0%)
All	All	0.65	0/7681	0.81	4/10382 (0.0%)

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 (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	459	ASN	N-CA-C	-5.53	104.28	111.74
1	A	457	ASP	N-CA-C	5.48	116.95	110.97
1	B	352	GLY	N-CA-C	-5.26	108.06	115.32
1	A	116	GLU	CB-CA-C	-5.19	110.57	116.54

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	244	LYS	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	3746	0	3627	113	0
1	B	3764	0	3648	71	0
2	A	3	0	0	2	0
2	B	4	0	0	1	0
3	A	5	0	0	1	0
3	B	13	0	0	2	0
4	A	18	0	24	2	0
4	B	24	0	32	2	0
5	A	199	0	0	6	0
5	B	200	0	0	6	0
All	All	7976	0	7331	179	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 12.

All (179) 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:A:358:ALA:CB	1:A:360:ASN:N	2.08	1.17
1:A:358:ALA:HB1	1:A:360:ASN:N	1.59	1.16
1:A:358:ALA:HB3	1:A:361:GLY:H	0.97	1.13
1:A:359:LYS:HG2	1:A:360:ASN:N	1.61	1.09
1:A:358:ALA:HB1	1:A:359:LYS:CB	1.81	1.09
1:A:359:LYS:CG	1:A:360:ASN:H	1.65	1.05
1:A:358:ALA:CB	1:A:359:LYS:C	2.29	1.05
1:B:457:ASP:OD1	1:B:458:TYR:N	1.94	1.01
1:A:358:ALA:HB3	1:A:361:GLY:N	1.76	0.99
1:A:358:ALA:HB1	1:A:359:LYS:C	1.86	0.99
1:B:158:ARG:HB3	1:B:159:PRO:HD3	1.43	0.97
1:A:359:LYS:HG2	1:A:360:ASN:H	0.81	0.95
1:A:301:ASP:HB3	1:B:431:LYS:HE2	1.50	0.94
1:A:358:ALA:HB1	1:A:360:ASN:H	1.31	0.91
1:B:158:ARG:HB3	1:B:159:PRO:CD	2.01	0.90
1:B:303:LYS:HZ1	1:B:456:CYS:HB3	1.36	0.90
1:A:324:GLN:HG3	1:A:326:THR:OG1	1.74	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:358:ALA:HB3	1:A:359:LYS:C	2.00	0.86
1:A:358:ALA:HB1	1:A:359:LYS:CA	2.06	0.86
1:A:353:LYS:HE2	1:A:368:GLU:CG	2.06	0.86
1:A:358:ALA:HB3	1:A:360:ASN:N	1.91	0.85
1:A:339:GLU:HG2	1:A:346:THR:OG1	1.76	0.85
1:A:358:ALA:HB1	1:A:359:LYS:CG	2.05	0.85
1:A:353:LYS:HE2	1:A:368:GLU:HG2	1.56	0.84
1:A:30:GLY:O	1:A:32:LYS:HG3	1.78	0.84
1:A:359:LYS:CG	1:A:360:ASN:N	2.30	0.83
1:A:358:ALA:HB1	1:A:359:LYS:HB3	1.63	0.81
1:A:30:GLY:C	1:A:32:LYS:HG3	2.06	0.81
1:A:359:LYS:O	1:A:361:GLY:N	2.16	0.79
1:A:305[B]:CYS:SG	1:A:338:ILE:HD11	2.21	0.78
1:B:303:LYS:NZ	1:B:457:ASP:H	1.81	0.78
1:A:359:LYS:O	1:A:360:ASN:C	2.26	0.77
1:A:358:ALA:HB1	1:A:359:LYS:HG2	1.64	0.77
1:A:218:SER:HB3	4:A:511:GOL:O2	1.84	0.77
1:B:158:ARG:CB	1:B:159:PRO:HD3	2.16	0.75
1:A:303:LYS:HE3	1:B:400:VAL:O	1.86	0.74
1:A:353:LYS:CE	1:A:368:GLU:HG2	2.18	0.73
1:A:281:LEU:HD21	1:A:325:SER:HB3	1.73	0.69
1:A:358:ALA:CB	1:A:359:LYS:CB	2.66	0.69
1:A:317:LEU:HB3	1:A:333:SER:HB2	1.75	0.68
1:A:486:ASP:HB2	1:A:487:PRO:CD	2.24	0.68
1:A:358:ALA:CB	1:A:359:LYS:HB3	2.24	0.67
1:B:284:ASN:OD1	1:B:285:GLN:HG2	1.96	0.66
1:B:303:LYS:HZ1	1:B:456:CYS:CB	2.09	0.66
1:B:28:ALA:HB2	5:B:698:HOH:O	1.96	0.65
1:B:351:ASN:HB3	1:B:353:LYS:H	1.62	0.64
1:A:358:ALA:CA	1:A:359:LYS:HB3	2.27	0.64
1:A:314:TYR:CZ	1:A:330:LYS:HE2	2.33	0.64
1:A:82:ARG:HG2	5:A:772:HOH:O	1.98	0.64
1:A:358:ALA:CB	1:A:361:GLY:H	1.92	0.63
1:A:217:ARG:NH1	1:A:248:VAL:HB	2.15	0.62
1:A:484:THR:HG21	1:B:477:VAL:CG2	2.30	0.62
1:A:339:GLU:OE2	1:A:346:THR:CG2	2.47	0.62
1:B:458:TYR:HB2	1:B:460:LYS:HE2	1.82	0.61
1:A:431:LYS:HE2	1:A:443:SER:O	2.01	0.60
1:A:88:ASP:HB2	2:A:502:BR:BR	2.57	0.59
1:B:303:LYS:NZ	1:B:456:CYS:HB3	2.15	0.59
1:B:88:ASP:HB2	2:B:502:BR:BR	2.58	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:229:ARG:HA	1:A:245:ALA:O	2.04	0.58
1:B:27:GLU:HB2	1:B:32:LYS:HG2	1.84	0.57
1:B:303:LYS:NZ	1:B:457:ASP:N	2.51	0.57
1:A:440:ALA:HA	1:A:479:LYS:HG2	1.86	0.57
1:A:486:ASP:HB2	1:A:487:PRO:HD3	1.85	0.57
1:A:353:LYS:HE2	1:A:368:GLU:CD	2.30	0.57
1:A:358:ALA:CB	1:A:360:ASN:H	1.98	0.57
1:A:317:LEU:CD2	1:A:351:ASN:HB2	2.35	0.56
1:B:351:ASN:ND2	1:B:353:LYS:HD3	2.20	0.56
1:A:377:LEU:HD12	5:A:688:HOH:O	2.05	0.55
1:B:343:ARG:HG2	1:B:452:PHE:HE1	1.71	0.55
1:B:457:ASP:CG	1:B:458:TYR:N	2.64	0.55
1:A:299:ASP:HB2	1:A:335:TYR:CE1	2.41	0.55
1:A:353:LYS:CD	1:A:368:GLU:HG2	2.37	0.55
1:A:302:THR:HA	1:B:444:SER:HB2	1.87	0.55
1:A:359:LYS:C	1:A:361:GLY:N	2.62	0.55
1:A:127:SER:HB2	1:A:128:PRO:HD2	1.89	0.54
1:A:362:GLN:HG3	1:A:363:LEU:N	2.21	0.54
1:B:258:GLN:NE2	5:B:796:HOH:O	2.29	0.54
1:A:304:LYS:HG2	1:A:337:ASP:OD1	2.07	0.54
5:A:729:HOH:O	1:B:400:VAL:HG13	2.07	0.54
1:B:183:ASP:O	1:B:184:GLN:CB	2.56	0.54
1:B:303:LYS:HZ3	1:B:457:ASP:H	1.56	0.54
1:B:264:VAL:HG21	1:B:288:GLU:HG2	1.90	0.54
1:A:339:GLU:OE2	1:A:346:THR:HG21	2.08	0.54
1:A:30:GLY:CA	1:A:32:LYS:HE2	2.38	0.53
1:A:330:LYS:HE3	3:A:505:CL:CL	2.45	0.53
1:A:351:ASN:OD1	1:A:353:LYS:HB2	2.08	0.53
1:B:100:ARG:HD2	1:B:184:GLN:NE2	2.23	0.53
1:A:358:ALA:HA	1:A:359:LYS:HB3	1.90	0.52
1:A:358:ALA:HB2	5:A:785:HOH:O	2.10	0.52
1:A:99:GLY:HA2	1:A:213:THR:OG1	2.10	0.51
1:B:82:ARG:HD3	3:B:507:CL:CL	2.47	0.51
1:A:27:GLU:OE1	1:A:32:LYS:HB2	2.10	0.50
1:A:317:LEU:HD21	1:A:351:ASN:HB2	1.92	0.50
1:A:156:SER:O	1:A:161:ASP:HA	2.12	0.50
1:B:269:ASN:HD21	1:B:283:ALA:HB3	1.77	0.49
1:A:339:GLU:OE2	1:A:341:ARG:HD2	2.13	0.49
1:B:354:PHE:O	1:B:365:ALA:HA	2.13	0.49
1:A:86:GLY:HA3	2:A:502:BR:BR	2.67	0.49
1:A:227:GLU:OE1	1:A:229:ARG:CZ	2.61	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:353:LYS:HE2	1:B:368:GLU:OE2	2.13	0.48
1:A:271:ARG:HD2	1:A:286:ASP:OD1	2.12	0.48
1:A:339:GLU:CG	1:A:346:THR:OG1	2.56	0.48
1:A:281:LEU:CD2	1:A:325:SER:HB3	2.41	0.48
1:B:300:ARG:HD2	1:B:458:TYR:OH	2.14	0.47
1:A:331:ASN:O	1:A:334:CYS:HB2	2.14	0.47
1:B:471:LYS:HD3	1:B:472:GLY:O	2.14	0.47
1:A:339:GLU:CD	1:A:341:ARG:HD2	2.39	0.47
1:A:398:ARG:HD2	1:A:403:THR:HB	1.96	0.47
1:B:304:LYS:NZ	1:B:337:ASP:OD1	2.41	0.47
1:A:490:LEU:HD13	1:B:400:VAL:HG11	1.97	0.47
1:B:82:ARG:CD	3:B:507:CL:CL	3.00	0.47
1:B:421:GLY:O	4:B:521:GOL:H32	2.15	0.46
1:B:246:THR:HB	5:B:619:HOH:O	2.15	0.46
1:A:427:ASP:OD1	1:A:431:LYS:HG2	2.16	0.46
1:A:471:LYS:HG2	1:A:482:ALA:HB2	1.97	0.46
1:A:30:GLY:HA3	1:A:32:LYS:HE2	1.97	0.46
1:B:406:ALA:HB1	1:B:491:TRP:HH2	1.80	0.46
1:B:458:TYR:CB	1:B:460:LYS:HE2	2.46	0.46
1:A:331:ASN:H	1:A:334:CYS:HB2	1.79	0.45
1:A:460:LYS:HE2	1:A:492:GLU:HG3	1.97	0.45
1:A:415:GLN:HB2	1:A:426:LYS:HG3	1.97	0.45
1:A:486:ASP:CB	1:A:487:PRO:CD	2.94	0.45
1:A:305[B]:CYS:SG	1:A:338:ILE:CD1	3.01	0.45
1:B:193:HIS:CD2	1:B:193:HIS:N	2.84	0.45
1:B:299:ASP:O	1:B:303:LYS:N	2.44	0.45
1:B:373:SER:OG	1:B:374:GLU:OE2	2.35	0.45
1:A:48:LEU:HD11	1:A:60:VAL:HG21	1.99	0.45
1:A:108:HIS:O	1:A:109:ARG:HB2	2.16	0.45
1:B:90:ARG:HD3	1:B:106:GLU:OE1	2.16	0.45
1:B:303:LYS:HZ2	1:B:457:ASP:N	2.13	0.45
1:B:42:LYS:HG2	5:B:666:HOH:O	2.17	0.45
1:B:183:ASP:O	1:B:184:GLN:HB3	2.17	0.45
1:B:447:THR:HB	1:B:448:PRO:HD2	1.98	0.45
1:B:421:GLY:O	4:B:521:GOL:C3	2.65	0.44
1:B:474:HIS:HD2	5:B:715:HOH:O	1.99	0.44
1:B:33:VAL:HG12	1:B:119:LEU:HD12	1.99	0.44
1:B:51:PRO:CB	1:B:52:PRO:HD2	2.48	0.44
1:B:260[A]:CYS:SG	1:B:296:LEU:O	2.76	0.44
1:A:353:LYS:HE2	1:A:368:GLU:OE2	2.18	0.44
1:A:161:ASP:OD2	5:A:797:HOH:O	2.21	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:339:GLU:CD	1:A:346:THR:HG1	2.26	0.43
1:B:379:LYS:HD3	1:B:418:PHE:CD2	2.54	0.43
1:B:300:ARG:CG	1:B:301:ASP:N	2.82	0.43
1:B:367:VAL:CG1	1:B:368:GLU:N	2.82	0.43
1:A:94:VAL:HB	1:A:102:SER:HB2	2.01	0.42
1:A:161:ASP:OD1	1:A:161:ASP:N	2.48	0.42
1:B:100:ARG:HD2	1:B:184:GLN:HE21	1.84	0.42
1:A:148:THR:HG21	1:A:251:ASP:OD2	2.19	0.42
1:A:182:GLN:OE1	1:A:187:SER:OG	2.32	0.42
1:A:217:ARG:HB2	4:A:511:GOL:H11	2.02	0.42
1:B:269:ASN:O	1:B:270:GLU:HB2	2.20	0.42
1:A:110:ARG:NH1	1:A:122:PHE:O	2.50	0.42
1:B:487:PRO:HD3	5:B:756:HOH:O	2.18	0.42
1:A:31:PHE:CZ	1:A:68:ARG:NH1	2.88	0.42
1:A:316:THR:O	1:A:323:VAL:HA	2.19	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	0.42
1:A:332:ALA:O	1:A:350:SER:HB2	2.20	0.42
1:B:434:THR:HG22	1:B:449:VAL:HG23	2.01	0.42
1:A:30:GLY:HA3	1:A:32:LYS:CD	2.50	0.41
1:B:10:VAL:HG21	1:B:142:VAL:HG12	2.00	0.41
1:B:353:LYS:CG	1:B:368:GLU:OE2	2.68	0.41
1:B:299:ASP:O	1:B:303:LYS:HA	2.20	0.41
1:A:97:ASP:O	1:A:185:ARG:NH2	2.45	0.41
1:B:282:SER:OG	1:B:284:ASN:ND2	2.54	0.41
1:A:78:VAL:HG23	1:A:120:SER:HA	2.02	0.41
1:A:447:THR:HA	1:A:448:PRO:HD3	1.93	0.41
1:B:40:LEU:HD22	1:B:135:HIS:CG	2.56	0.41
1:A:55:ALA:HA	1:A:56:GLY:HA2	1.75	0.41
1:A:356:THR:OG1	1:A:357:SER:N	2.54	0.41
1:B:99:GLY:HA2	1:B:213:THR:OG1	2.21	0.41
1:B:405:ASP:HB2	1:B:408:ARG:CD	2.51	0.41
1:A:90:ARG:HD3	1:A:106:GLU:OE1	2.20	0.40
1:B:73:ASP:OD1	1:B:73:ASP:C	2.64	0.40
1:A:401:THR:CB	1:A:403:THR:HG1	2.34	0.40
1:B:432:TYR:CD2	1:B:448:PRO:HB3	2.56	0.40
1:A:295:GLN:HB3	1:A:308:ARG:HB3	2.04	0.40
1:A:298:ILE:HB	1:A:457:ASP:HA	2.02	0.40
1:A:399:LYS:HA	5:A:667:HOH:O	2.20	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	474/493 (96%)	447 (94%)	27 (6%)	0	100	100
1	B	478/493 (97%)	456 (95%)	22 (5%)	0	100	100
All	All	952/986 (97%)	903 (95%)	49 (5%)	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	395/403 (98%)	391 (99%)	4 (1%)	68	77
1	B	398/403 (99%)	396 (100%)	2 (0%)	81	87
All	All	793/806 (98%)	787 (99%)	6 (1%)	73	80

All (6) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	84	VAL
1	A	138	MET
1	A	287	GLU
1	A	323	VAL
1	B	138	MET
1	B	193	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11)

such sidechains are listed below:

Mol	Chain	Res	Type
1	A	65	HIS
1	A	193	HIS
1	A	198	HIS
1	A	258	GLN
1	A	285	GLN
1	A	310	HIS
1	A	459	ASN
1	B	124	GLN
1	B	277	GLN
1	B	331	ASN
1	B	459	ASN

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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 32 ligands modelled in this entry, 25 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	519	-	5,5,5	0.28	0	5,5,5	0.28	0
4	GOL	A	510	-	5,5,5	0.39	0	5,5,5	0.73	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	GOL	B	521	-	5,5,5	0.47	0	5,5,5	0.77	0
4	GOL	B	518	-	5,5,5	0.17	0	5,5,5	0.71	0
4	GOL	A	511	-	5,5,5	0.28	0	5,5,5	0.25	0
4	GOL	A	509	-	5,5,5	0.28	0	5,5,5	0.44	0
4	GOL	B	520	-	5,5,5	0.33	0	5,5,5	0.19	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	519	-	-	2/4/4/4	-
4	GOL	A	510	-	-	4/4/4/4	-
4	GOL	B	521	-	-	2/4/4/4	-
4	GOL	B	518	-	-	2/4/4/4	-
4	GOL	A	511	-	-	4/4/4/4	-
4	GOL	A	509	-	-	2/4/4/4	-
4	GOL	B	520	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	510	GOL	O1-C1-C2-C3
4	A	511	GOL	C1-C2-C3-O3
4	B	519	GOL	C1-C2-C3-O3
4	B	519	GOL	O2-C2-C3-O3
4	A	511	GOL	O2-C2-C3-O3
4	B	521	GOL	O2-C2-C3-O3
4	A	509	GOL	O1-C1-C2-C3
4	A	510	GOL	C1-C2-C3-O3
4	A	511	GOL	O1-C1-C2-C3
4	B	518	GOL	O1-C1-C2-C3
4	B	520	GOL	C1-C2-C3-O3
4	B	521	GOL	C1-C2-C3-O3
4	A	510	GOL	O1-C1-C2-O2

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Mol	Chain	Res	Type	Atoms
4	A	510	GOL	O2-C2-C3-O3
4	A	511	GOL	O1-C1-C2-O2
4	A	509	GOL	O1-C1-C2-O2
4	B	518	GOL	O1-C1-C2-O2
4	B	520	GOL	O1-C1-C2-O2
4	B	520	GOL	O2-C2-C3-O3

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	521	GOL	2	0
4	A	511	GOL	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	480/493 (97%)	0.25	22 (4%) 37 36	24, 44, 69, 89	2 (0%)
1	B	481/493 (97%)	0.24	21 (4%) 39 38	24, 42, 90, 122	3 (0%)
All	All	961/986 (97%)	0.24	43 (4%) 38 37	24, 43, 79, 122	5 (0%)

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	400	VAL	3.9
1	A	332	ALA	3.9
1	A	29	PHE	3.7
1	A	370	ALA	3.6
1	A	55	ALA	3.5
1	B	55	ALA	3.5
1	A	7	ALA	3.4
1	B	343	ARG	3.3
1	A	358	ALA	3.1
1	A	218	SER	3.0
1	B	351	ASN	3.0
1	A	157	ALA	2.9
1	A	160	ALA	2.8
1	B	457	ASP	2.7
1	B	361	GLY	2.7
1	B	327	ALA	2.6
1	B	318	THR	2.5
1	B	328	SER	2.5
1	A	245	ALA	2.5
1	B	326	THR	2.4
1	B	357	SER	2.4
1	B	366	SER	2.4
1	B	399	LYS	2.4
1	A	248	VAL	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	29	PHE	2.4
1	B	56	GLY	2.3
1	A	51	PRO	2.3
1	A	444	SER	2.3
1	A	457	ASP	2.3
1	A	219	GLY	2.3
1	B	338	ILE	2.3
1	B	317	LEU	2.3
1	B	286	ASP	2.3
1	A	317	LEU	2.2
1	B	345	ILE	2.1
1	B	336	PHE	2.1
1	B	365	ALA	2.1
1	A	401	THR	2.0
1	A	301	ASP	2.0
1	B	314	TYR	2.0
1	A	300	ARG	2.0
1	A	246	THR	2.0
1	A	319	ALA	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
4	GOL	B	521	6/6	0.75	0.18	47,51,54,57	0
4	GOL	B	520	6/6	0.81	0.15	58,60,62,65	0
4	GOL	B	519	6/6	0.82	0.13	54,58,60,63	0
4	GOL	A	511	6/6	0.85	0.12	48,53,54,57	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	CL	B	515	1/1	0.85	0.14	66,66,66,66	0
4	GOL	A	509	6/6	0.88	0.12	35,43,44,46	0
3	CL	B	516	1/1	0.89	0.18	82,82,82,82	0
3	CL	A	505	1/1	0.89	0.10	66,66,66,66	0
3	CL	B	514	1/1	0.90	0.11	66,66,66,66	0
3	CL	A	508	1/1	0.90	0.10	62,62,62,62	0
4	GOL	A	510	6/6	0.91	0.10	38,44,46,46	0
4	GOL	B	518	6/6	0.92	0.09	38,41,42,42	0
3	CL	A	506	1/1	0.92	0.09	62,62,62,62	0
3	CL	B	510	1/1	0.94	0.08	53,53,53,53	0
3	CL	B	512	1/1	0.94	0.08	58,58,58,58	0
3	CL	A	507	1/1	0.94	0.08	55,55,55,55	0
3	CL	B	508	1/1	0.94	0.11	44,44,44,44	0
2	BR	A	502	1/1	0.95	0.08	56,56,56,56	0
3	CL	B	517	1/1	0.95	0.06	51,51,51,51	0
3	CL	B	513	1/1	0.95	0.11	68,68,68,68	0
2	BR	A	503	1/1	0.96	0.10	65,65,65,65	0
3	CL	A	504	1/1	0.96	0.07	57,57,57,57	0
3	CL	B	507	1/1	0.96	0.09	37,37,37,37	0
2	BR	B	504	1/1	0.97	0.07	66,66,66,66	0
3	CL	B	509	1/1	0.97	0.08	45,45,45,45	0
3	CL	B	506	1/1	0.98	0.09	41,41,41,41	0
2	BR	A	501	1/1	0.98	0.06	45,45,45,45	0
3	CL	B	511	1/1	0.98	0.05	57,57,57,57	0
3	CL	B	505	1/1	0.98	0.05	34,34,34,34	0
2	BR	B	501	1/1	0.99	0.04	44,44,44,44	0
2	BR	B	502	1/1	0.99	0.04	46,46,46,46	0
2	BR	B	503	1/1	1.00	0.05	43,43,43,43	0

6.5 Other polymers

There are no such residues in this entry.