Full wwPDB X-ray Structure Validation Report (i)

Nov 15, 2023 – 02:24 PM JST

PDB ID	:	6JEG
Title	:	Crystal structure of calcium free human gelsolin amyloid mutant G167R
Authors	:	Zorgati, H.; Robinson, R.C.
Deposited on	:	2019-02-05
Resolution	:	2.98 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 (i) 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 (1)) were used in the production of this report:

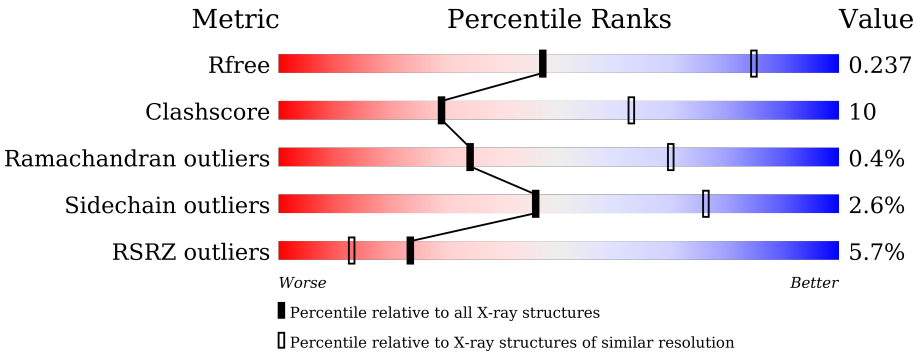
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.98 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	2754 (3.00-2.96)
Clashscore	141614	3103 (3.00-2.96)
Ramachandran outliers	138981	2993 (3.00-2.96)
Sidechain outliers	138945	2996 (3.00-2.96)
RSRZ outliers	127900	2644 (3.00-2.96)

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 <=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	729	5% 78% 19%
1	B	729	7% 80% 17%

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11203 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.

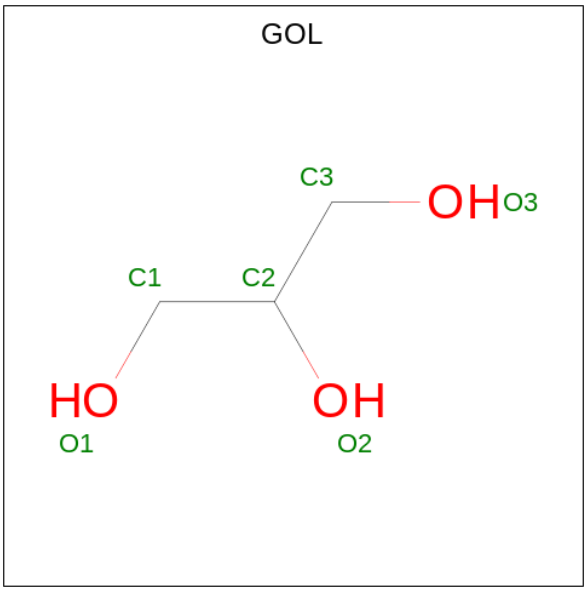
- Molecule 1 is a protein called Gelsolin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	714	Total	C	N	O	S	0	0	0
			5580	3523	978	1063	16			
1	B	718	Total	C	N	O	S	0	0	0
			5608	3538	982	1072	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	167	ARG	GLY	engineered mutation	UNP P06396
B	167	ARG	GLY	engineered mutation	UNP P06396

- Molecule 2 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

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

- Molecule 3 is water.

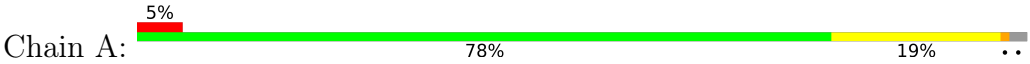
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	3	Total O 3 3	0	0

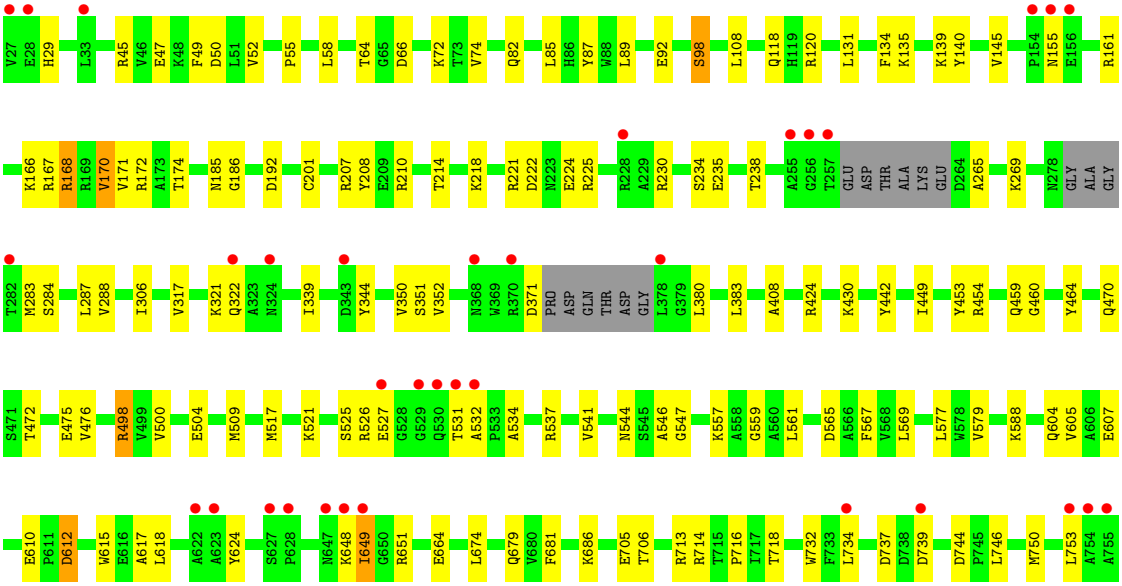
3 Residue-property plots (i)

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: Gelsolin

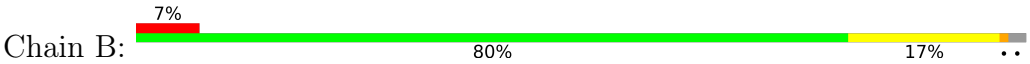
Chain A:

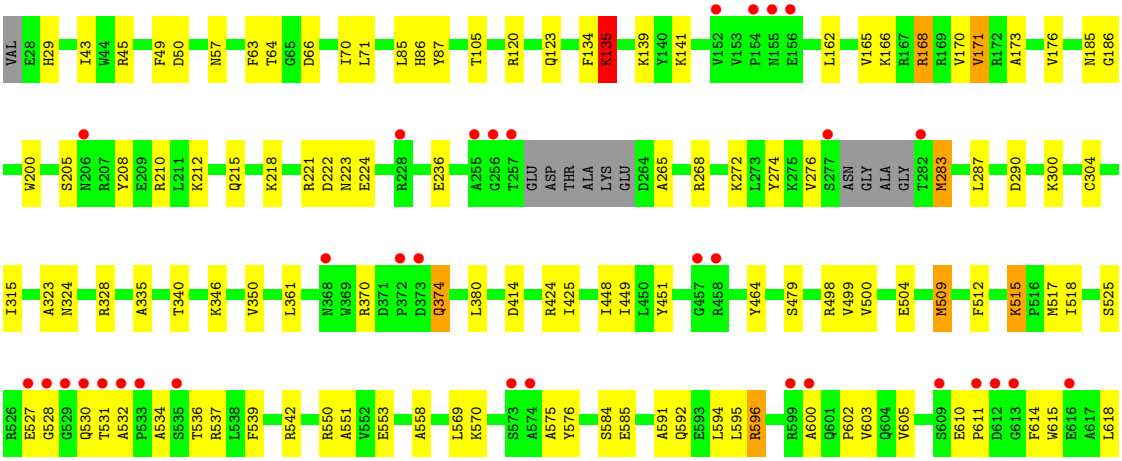


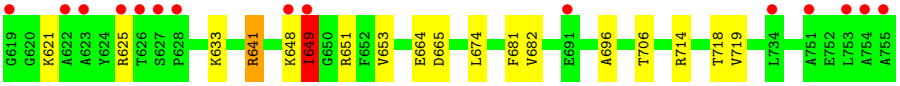


• Molecule 1: Gelsolin

Chain B:






G619
G620
K621
A622
A623
Y624
R625
T626
S627
P628
K633
R641
K648
I649
G650
R651
F652
V653
E664
D665
L674
F681
V682
E691
A696
T706
R714
T718
V719
L734
A751
E752
L753
A754
A755

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α, β, γ	172.27Å 172.27Å 150.11Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	19.78 – 2.98 19.78 – 2.97	Depositor EDS
% Data completeness (in resolution range)	96.4 (19.78-2.98) 96.4 (19.78-2.97)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.59 (at 2.98Å)	Xtriage
Refinement program	PHENIX (1.13_2998: ???)	Depositor
R, R_{free}	0.198 , 0.237 0.198 , 0.237	Depositor DCC
R_{free} test set	2334 reflections (5.18%)	wwPDB-VP
Wilson B-factor (Å^2)	56.5	Xtriage
Anisotropy	0.019	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 40.0	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	11203	wwPDB-VP
Average B, all atoms (Å^2)	57.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.36% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ 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: GOL

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	0.33	0/5706	0.58	5/7726 (0.1%)
1	B	0.32	0/5736	0.56	5/7769 (0.1%)
All	All	0.32	0/11442	0.57	10/15495 (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	2
All	All	0	3

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	230	ARG	CG-CD-NE	-13.54	83.38	111.80
1	B	649	ILE	CG1-CB-CG2	-10.77	87.71	111.40
1	B	135	LYS	CD-CE-NZ	8.58	131.44	111.70
1	A	579	VAL	CG1-CB-CG2	-8.08	97.97	110.90
1	B	283	MET	CB-CG-SD	-7.77	89.09	112.40
1	A	588	LYS	CA-CB-CG	7.39	129.65	113.40
1	A	230	ARG	NE-CZ-NH2	-7.18	116.71	120.30
1	B	283	MET	CG-SD-CE	-5.43	91.51	100.20
1	B	135	LYS	CA-CB-CG	5.27	124.99	113.40
1	A	588	LYS	N-CA-CB	5.07	119.72	110.60

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	135	LYS	Peptide
1	B	135	LYS	Peptide
1	B	374	GLN	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	5580	0	5445	97	0
1	B	5608	0	5464	116	0
2	A	6	0	8	3	0
2	B	6	0	8	1	0
3	B	3	0	0	0	0
All	All	11203	0	10925	213	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 (213) 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:570:LYS:NZ	1:B:600:ALA:HB2	1.17	1.42
1:B:570:LYS:NZ	1:B:600:ALA:CB	1.93	1.31
1:B:592:GLN:HA	1:B:595:LEU:HD23	1.30	1.08
1:A:168:ARG:NH1	1:A:208:TYR:CD1	2.21	1.07
1:B:570:LYS:HZ1	1:B:600:ALA:CB	1.69	0.94
1:A:168:ARG:NH1	1:A:208:TYR:HD1	1.59	0.94
1:B:168:ARG:NH1	1:B:208:TYR:CD2	2.36	0.93
1:B:570:LYS:HZ2	1:B:600:ALA:HB2	1.11	0.92
1:B:570:LYS:HZ3	1:B:600:ALA:HB2	1.21	0.92
1:B:168:ARG:HH11	1:B:208:TYR:HD2	1.11	0.90
1:B:570:LYS:HZ2	1:B:600:ALA:CB	1.69	0.89
1:B:576:TYR:HE1	1:B:614:PHE:HA	1.39	0.88
1:A:454:ARG:HH11	1:A:459:GLN:HG3	1.42	0.84
1:A:118:GLN:HB2	1:A:352:VAL:HG22	1.63	0.80

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:682:VAL:HG11	1:B:696:ALA:HB1	1.67	0.77
1:A:139:LYS:NZ	1:A:224:GLU:OE1	2.23	0.72
1:B:168:ARG:NH1	1:B:208:TYR:HD2	1.80	0.71
1:B:527:GLU:O	1:B:527:GLU:HG2	1.91	0.69
1:B:139:LYS:NZ	1:B:224:GLU:OE1	2.25	0.69
1:B:648:LYS:HG3	1:B:649:ILE:HD13	1.72	0.69
1:A:55:PRO:HB2	1:A:58:LEU:HD12	1.74	0.69
1:B:592:GLN:HA	1:B:595:LEU:CD2	2.16	0.69
1:A:131:LEU:HD22	1:A:170:VAL:HG11	1.73	0.69
1:B:512:PHE:CE2	1:B:517:MET:HG3	2.29	0.68
1:B:570:LYS:HZ1	1:B:600:ALA:HB1	1.56	0.68
1:B:323:ALA:O	1:B:328:ARG:NH1	2.27	0.67
1:A:648:LYS:HG3	1:A:649:ILE:HD13	1.77	0.67
1:B:570:LYS:NZ	1:B:600:ALA:HB1	2.03	0.67
1:B:591:ALA:O	1:B:595:LEU:HD22	1.96	0.66
1:A:207:ARG:NH2	1:A:744:ASP:OD2	2.30	0.64
1:A:45:ARG:NH1	1:A:66:ASP:OD2	2.30	0.64
1:B:615:TRP:CE3	1:B:621:LYS:HD2	2.33	0.64
1:A:339:ILE:HG23	1:A:344:TYR:HB2	1.81	0.63
1:B:425:ILE:HD13	1:B:479:SER:HA	1.82	0.62
1:A:168:ARG:HD2	1:A:208:TYR:HD1	1.65	0.62
1:B:576:TYR:CE1	1:B:614:PHE:HA	2.30	0.62
1:A:504:GLU:HG2	1:A:517:MET:HE2	1.82	0.61
1:A:679:GLN:OE1	2:A:801:GOL:H12	2.01	0.60
1:B:212:LYS:NZ	1:B:665:ASP:OD1	2.34	0.60
1:B:576:TYR:HA	1:B:603:VAL:HG22	1.81	0.60
1:A:140:TYR:HD1	1:A:174:THR:HG22	1.67	0.60
1:A:234:SER:HB2	1:A:238:THR:HG23	1.84	0.59
1:A:265:ALA:HB3	1:A:651:ARG:HH21	1.66	0.59
1:A:737:ASP:OD1	1:A:739:ASP:N	2.35	0.59
1:A:664:GLU:OE2	1:A:664:GLU:N	2.35	0.59
1:B:605:VAL:HG13	1:B:610:GLU:HB3	1.85	0.59
1:B:682:VAL:HB	1:B:719:VAL:HA	1.84	0.59
1:A:521:LYS:O	1:A:521:LYS:NZ	2.26	0.58
1:B:43:ILE:HD13	1:B:70:ILE:CG1	2.32	0.58
1:B:168:ARG:NH1	1:B:208:TYR:HB3	2.18	0.58
1:A:321:LYS:C	1:A:322:GLN:HG2	2.24	0.58
1:B:361:LEU:HD13	1:B:641:ARG:NH2	2.19	0.57
1:A:89:LEU:HD22	1:A:98:SER:HB2	1.87	0.57
1:B:414:ASP:OD2	1:B:451:TYR:OH	2.18	0.56
1:A:168:ARG:HH11	1:A:208:TYR:HD1	0.75	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:43:ILE:HD13	1:B:70:ILE:HG13	1.86	0.56
1:B:64:THR:HG23	1:B:123:GLN:OE1	2.06	0.56
1:B:576:TYR:HE1	1:B:614:PHE:CA	2.13	0.56
1:B:584:SER:OG	1:B:585:GLU:OE1	2.24	0.55
1:A:168:ARG:HD2	1:A:208:TYR:CD1	2.40	0.55
1:A:470:GLN:HG3	1:A:521:LYS:HZ1	1.71	0.55
1:A:269:LYS:HG2	1:A:706:THR:HA	1.89	0.55
1:B:527:GLU:O	1:B:530:GLN:NE2	2.40	0.54
1:B:531:THR:HG22	1:B:532:ALA:H	1.73	0.54
1:A:72:LYS:HE3	1:A:74:VAL:HG12	1.90	0.54
1:B:509:MET:HE2	1:B:550:ARG:NE	2.23	0.54
1:A:504:GLU:HG2	1:A:517:MET:CE	2.38	0.54
1:B:276:VAL:HG13	1:B:283:MET:CE	2.38	0.54
1:B:361:LEU:CD1	1:B:641:ARG:CZ	2.85	0.53
1:B:500:VAL:HB	1:B:714:ARG:HB3	1.88	0.53
1:B:45:ARG:NH1	1:B:66:ASP:OD2	2.41	0.53
1:A:498:ARG:HH12	1:A:716:PRO:HA	1.73	0.53
1:A:306:ILE:HG12	1:A:317:VAL:HG12	1.91	0.53
1:B:596:ARG:O	1:B:596:ARG:NE	2.39	0.53
1:B:674:LEU:HB3	1:B:681:PHE:HB2	1.91	0.53
1:A:454:ARG:HD2	1:A:459:GLN:HG3	1.90	0.52
1:A:207:ARG:NH1	1:A:746:LEU:HD23	2.24	0.52
1:A:577:LEU:HD22	1:A:604:GLN:HG3	1.92	0.52
1:B:539:PHE:HE1	1:B:553:GLU:HB2	1.75	0.52
1:A:544:ASN:OD1	1:A:547:GLY:N	2.42	0.51
1:A:569:LEU:HD22	1:A:618:LEU:HD21	1.92	0.51
1:B:272:LYS:HG2	1:B:290:ASP:O	2.10	0.51
1:A:161:ARG:HH21	1:A:225:ARG:HH22	1.58	0.51
1:B:200:TRP:CZ2	1:B:236:GLU:HG2	2.45	0.51
1:B:648:LYS:HG3	1:B:649:ILE:CD1	2.40	0.51
1:A:120:ARG:HB3	1:A:350:VAL:HB	1.93	0.51
1:A:287:LEU:HD12	1:A:288:VAL:H	1.76	0.51
1:B:534:ALA:HB3	1:B:537:ARG:HG2	1.93	0.50
1:B:374:GLN:HB2	1:B:380:LEU:HD23	1.92	0.50
1:B:539:PHE:CD2	1:B:594:LEU:HD21	2.47	0.50
1:B:361:LEU:HD13	1:B:641:ARG:CZ	2.41	0.50
1:A:185:ASN:OD1	1:A:186:GLY:N	2.44	0.50
1:A:531:THR:HG22	1:A:532:ALA:H	1.77	0.50
1:B:315:ILE:HD11	1:B:335:ALA:HB1	1.94	0.50
1:A:454:ARG:HH11	1:A:459:GLN:CG	2.18	0.49
1:A:214:THR:O	1:A:218:LYS:HG3	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:557:LYS:HG2	1:A:559:GLY:H	1.76	0.49
1:B:499:VAL:HG21	1:B:504:GLU:HA	1.94	0.49
1:B:539:PHE:CE1	1:B:553:GLU:HB2	2.47	0.49
1:A:534:ALA:HB3	1:A:537:ARG:HG2	1.95	0.49
1:A:168:ARG:NH1	1:A:208:TYR:HB3	2.28	0.49
1:A:472:THR:O	1:A:476:VAL:HG23	2.13	0.49
1:A:500:VAL:HB	1:A:714:ARG:HB3	1.94	0.49
1:B:504:GLU:HG2	1:B:517:MET:CE	2.43	0.49
1:B:425:ILE:HD13	1:B:479:SER:CA	2.43	0.48
1:B:424:ARG:NH1	1:B:525:SER:OG	2.45	0.48
1:B:185:ASN:OD1	1:B:186:GLY:N	2.46	0.48
1:B:509:MET:CE	1:B:550:ARG:HE	2.26	0.48
1:B:70:ILE:HB	1:B:86:HIS:HB2	1.95	0.48
1:B:218:LYS:HE3	1:B:222:ASP:OD2	2.13	0.48
1:A:29:HIS:NE2	1:A:50:ASP:OD1	2.41	0.48
1:B:649:ILE:HG22	1:B:649:ILE:O	2.13	0.48
1:A:453:TYR:CZ	1:A:460:GLY:HA3	2.49	0.48
1:A:674:LEU:HB3	1:A:681:PHE:HB2	1.96	0.48
1:B:265:ALA:HB2	1:B:653:VAL:HA	1.96	0.48
1:B:425:ILE:HD13	1:B:479:SER:CB	2.44	0.47
1:B:504:GLU:HG2	1:B:517:MET:HE3	1.96	0.47
1:B:205:SER:OG	1:B:210:ARG:NH1	2.48	0.47
1:B:218:LYS:HD3	1:B:221:ARG:NH2	2.29	0.47
1:A:161:ARG:HH21	1:A:225:ARG:NH2	2.13	0.47
1:A:557:LYS:HG2	1:A:559:GLY:N	2.30	0.47
1:B:29:HIS:NE2	1:B:50:ASP:OD1	2.45	0.47
1:A:66:ASP:OD1	1:A:145:VAL:HG23	2.15	0.47
1:A:214:THR:HG23	1:A:753:LEU:HD23	1.97	0.47
1:A:317:VAL:HG22	1:A:352:VAL:HG12	1.96	0.47
1:B:340:THR:HG22	1:B:346:LYS:NZ	2.30	0.47
1:A:87:TYR:OH	1:A:120:ARG:HD2	2.15	0.47
1:A:269:LYS:HE2	1:A:705:GLU:O	2.15	0.47
1:B:592:GLN:O	1:B:595:LEU:N	2.48	0.46
1:B:168:ARG:HH12	1:B:208:TYR:CB	2.29	0.46
1:B:374:GLN:HB3	1:B:633:LYS:NZ	2.31	0.46
1:B:71:LEU:HD13	1:B:85:LEU:HD13	1.97	0.46
1:A:544:ASN:HD21	1:A:546:ALA:HB3	1.79	0.45
1:B:509:MET:HE2	1:B:550:ARG:HE	1.80	0.45
1:B:569:LEU:CD1	1:B:618:LEU:HD11	2.47	0.45
1:A:218:LYS:HB3	1:A:221:ARG:NH2	2.30	0.45
1:B:139:LYS:HG2	1:B:173:ALA:HB3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:686:LYS:HD2	1:A:686:LYS:HA	1.74	0.45
1:B:425:ILE:HD11	1:B:448:ILE:HG13	1.98	0.45
1:A:201:CYS:HB3	1:A:210:ARG:HD3	1.99	0.45
1:A:408:ALA:HB1	1:A:732:TRP:CD1	2.52	0.45
1:B:268:ARG:NH1	1:B:706:THR:OG1	2.50	0.45
1:A:85:LEU:HD21	1:A:108:LEU:HD23	1.99	0.45
1:B:87:TYR:HB3	1:B:105:THR:HG21	1.98	0.45
1:B:168:ARG:HH12	1:B:208:TYR:HB3	1.80	0.45
1:A:544:ASN:ND2	1:A:546:ALA:HB3	2.31	0.45
1:A:605:VAL:CG1	1:A:610:GLU:HB3	2.47	0.44
1:A:449:ILE:HB	1:A:464:TYR:HB2	1.99	0.44
1:B:605:VAL:CG1	1:B:610:GLU:HB3	2.48	0.44
1:A:221:ARG:NH1	1:A:222:ASP:OD2	2.51	0.44
1:B:57:ASN:O	1:B:223:ASN:HB3	2.17	0.44
1:B:610:GLU:HG3	1:B:614:PHE:HD2	1.82	0.44
1:A:605:VAL:HG11	1:A:610:GLU:HB3	1.99	0.44
1:A:45:ARG:NE	1:A:47:GLU:OE2	2.50	0.44
1:B:120:ARG:HB3	1:B:350:VAL:HB	2.00	0.44
1:A:82:GLN:HE22	1:A:383:LEU:H	1.65	0.44
1:B:504:GLU:OE1	1:B:550:ARG:NH1	2.49	0.44
1:B:218:LYS:NZ	1:B:221:ARG:HE	2.15	0.44
1:B:536:THR:HG23	1:B:569:LEU:HD21	1.99	0.44
1:A:475:GLU:OE1	1:A:526:ARG:NH1	2.49	0.43
1:B:542:ARG:HD2	1:B:550:ARG:HH21	1.83	0.43
1:B:570:LYS:HE3	1:B:575:ALA:HB2	2.01	0.43
1:A:454:ARG:NH1	1:A:459:GLN:OE1	2.52	0.43
1:A:648:LYS:O	1:A:649:ILE:HB	2.17	0.43
1:A:559:GLY:HA2	1:A:624:TYR:HB3	2.01	0.43
1:A:679:GLN:OE1	2:A:801:GOL:C1	2.65	0.43
1:A:118:GLN:O	1:A:351:SER:HA	2.19	0.42
1:B:576:TYR:N	1:B:576:TYR:CD2	2.85	0.42
1:B:276:VAL:HG13	1:B:283:MET:HE3	2.00	0.42
1:B:87:TYR:OH	1:B:120:ARG:HD2	2.19	0.42
1:A:192:ASP:CG	1:A:225:ARG:HH12	2.22	0.42
1:A:612:ASP:OD1	1:A:612:ASP:N	2.52	0.42
1:B:63:PHE:CZ	1:B:141:LYS:HE3	2.55	0.42
1:A:442:TYR:CD1	1:A:525:SER:HB3	2.54	0.42
1:B:515:LYS:H	1:B:515:LYS:HD2	1.85	0.42
1:B:539:PHE:HD2	1:B:594:LEU:HD21	1.84	0.42
1:B:558:ALA:HA	1:B:618:LEU:HD23	2.02	0.42
1:B:595:LEU:HG	1:B:602:PRO:HB3	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:ARG:NH1	1:B:208:TYR:CG	2.85	0.42
1:A:85:LEU:HD23	1:A:85:LEU:HA	1.86	0.42
1:A:610:GLU:OE2	1:A:615:TRP:CZ2	2.73	0.42
1:A:750:MET:HA	1:A:753:LEU:HG	2.02	0.42
1:B:215:GLN:NE2	1:B:664:GLU:OE1	2.52	0.42
1:B:718:THR:HG21	2:B:801:GOL:H32	2.02	0.42
1:A:561:LEU:HA	1:A:567:PHE:CZ	2.55	0.41
1:B:449:ILE:HB	1:B:464:TYR:HB2	2.02	0.41
1:B:162:LEU:HG	1:B:176:VAL:HG23	2.00	0.41
1:A:607:GLU:OE2	1:A:624:TYR:OH	2.33	0.41
1:A:66:ASP:CG	1:A:145:VAL:HG23	2.41	0.41
1:A:569:LEU:HD21	1:A:617:ALA:HB3	2.03	0.41
1:A:718:THR:HG21	2:A:801:GOL:H32	2.01	0.41
1:B:610:GLU:OE1	1:B:615:TRP:CZ2	2.74	0.41
1:A:424:ARG:NH1	1:A:525:SER:OG	2.52	0.41
1:A:651:ARG:HA	1:A:651:ARG:HD2	1.94	0.41
1:B:135:LYS:HB2	1:B:135:LYS:HE2	1.46	0.41
1:B:610:GLU:HB2	1:B:611:PRO:CD	2.51	0.41
1:B:265:ALA:O	1:B:268:ARG:HG2	2.21	0.41
1:B:274:TYR:CZ	1:B:287:LEU:HD13	2.55	0.41
1:B:276:VAL:CG1	1:B:283:MET:CE	2.99	0.41
1:B:518:ILE:HG23	1:B:551:ALA:HB3	2.02	0.41
1:B:165:VAL:HG12	1:B:171:VAL:HG22	2.02	0.41
1:A:235:GLU:H	1:A:238:THR:HG22	1.85	0.41
1:A:541:VAL:O	1:A:565:ASP:HB3	2.21	0.41
1:B:276:VAL:HB	1:B:304:CYS:O	2.20	0.41
1:A:47:GLU:CD	1:A:52:VAL:HG11	2.41	0.40
1:A:430:LYS:HB2	1:A:430:LYS:HE2	1.87	0.40
1:B:512:PHE:HE2	1:B:517:MET:HG3	1.85	0.40
1:B:218:LYS:HZ3	1:B:221:ARG:HH21	1.69	0.40
1:B:610:GLU:OE1	1:B:615:TRP:CH2	2.74	0.40
1:A:371:ASP:OD1	1:A:380:LEU:HD22	2.22	0.40
1:A:509:MET:CE	1:A:544:ASN:HB3	2.51	0.40
1:A:64:THR:HG22	1:A:92:GLU:H	1.87	0.40
1:A:168:ARG:NH1	1:A:208:TYR:CG	2.81	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	706/729 (97%)	673 (95%)	31 (4%)	2 (0%)	41	74
1	B	712/729 (98%)	678 (95%)	31 (4%)	3 (0%)	34	70
All	All	1418/1458 (97%)	1351 (95%)	62 (4%)	5 (0%)	34	70

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	649	ILE
1	B	528	GLY
1	A	170	VAL
1	B	170	VAL
1	B	649	ILE

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	584/594 (98%)	568 (97%)	16 (3%)	44	75
1	B	587/594 (99%)	572 (97%)	15 (3%)	46	76
All	All	1171/1188 (99%)	1140 (97%)	31 (3%)	46	76

All (31) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	PHE

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Mol	Chain	Res	Type
1	A	98	SER
1	A	134	PHE
1	A	155	ASN
1	A	166	LYS
1	A	167	ARG
1	A	168	ARG
1	A	171	VAL
1	A	172	ARG
1	A	283	MET
1	A	284	SER
1	A	498	ARG
1	A	527	GLU
1	A	612	ASP
1	A	713	ARG
1	A	734	LEU
1	B	49	PHE
1	B	134	PHE
1	B	166	LYS
1	B	168	ARG
1	B	171	VAL
1	B	300	LYS
1	B	324	ASN
1	B	370	ARG
1	B	498	ARG
1	B	509	MET
1	B	515	LYS
1	B	596	ARG
1	B	625	ARG
1	B	641	ARG
1	B	651	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	82	GLN
1	A	155	ASN
1	A	267	ASN
1	B	473	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

There are no monosaccharides in this entry.

5.6 Ligand geometry

2 ligands are modelled in this entry.

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	$\#\|Z\| > 2$
2	GOL	A	801	-	5,5,5	0.50	0	5,5,5	0.66	0
2	GOL	B	801	-	5,5,5	0.51	0	5,5,5	0.66	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
2	GOL	A	801	-	-	0/4/4/4	-
2	GOL	B	801	-	-	0/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	801	GOL	3	0
2	B	801	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, 95^{th} 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	714/729 (97%)	0.00	34 (4%)	30	18	22, 53, 98, 130	0
1	B	718/729 (98%)	0.07	48 (6%)	17	9	24, 52, 114, 154	0
All	All	1432/1458 (98%)	0.04	82 (5%)	23	13	22, 52, 104, 154	0

All (82) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	457	GLY	9.1
1	B	155	ASN	6.6
1	B	458	ARG	5.5
1	A	156	GLU	5.4
1	B	755	ALA	5.1
1	B	257	THR	4.8
1	B	372	PRO	4.6
1	B	282	THR	4.6
1	B	599	ARG	4.5
1	A	649	ILE	4.4
1	B	255	ALA	4.4
1	B	529	GLY	4.3
1	B	649	ILE	4.3
1	B	616	GLU	4.1
1	A	623	ALA	4.1
1	A	257	THR	4.1
1	B	619	GLY	4.0
1	B	611	PRO	4.0
1	A	754	ALA	3.9
1	A	755	ALA	3.9
1	A	622	ALA	3.8
1	B	373	ASP	3.7
1	A	648	LYS	3.7
1	B	256	GLY	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	612	ASP	3.6
1	B	530	GLN	3.6
1	A	256	GLY	3.5
1	B	531	THR	3.4
1	A	255	ALA	3.3
1	A	378	LEU	3.3
1	B	609	SER	3.2
1	B	277	SER	3.2
1	A	154	PRO	3.2
1	A	531	THR	3.2
1	B	206	ASN	3.1
1	A	282	THR	3.0
1	B	533	PRO	3.0
1	B	532	ALA	2.9
1	B	574	ALA	2.9
1	A	33	LEU	2.9
1	B	228	ARG	2.9
1	B	754	ALA	2.8
1	B	751	ALA	2.8
1	B	527	GLU	2.7
1	B	623	ALA	2.7
1	A	627	SER	2.7
1	B	600	ALA	2.7
1	B	626	THR	2.7
1	A	27	VAL	2.6
1	A	628	PRO	2.6
1	A	753	LEU	2.6
1	A	370	ARG	2.6
1	A	532	ALA	2.6
1	A	530	GLN	2.5
1	A	322	GLN	2.5
1	B	573	SER	2.5
1	A	739	ASP	2.4
1	B	625	ARG	2.4
1	A	527	GLU	2.4
1	B	368	ASN	2.4
1	B	528	GLY	2.4
1	A	529	GLY	2.3
1	A	324	ASN	2.3
1	B	156	GLU	2.3
1	B	648	LYS	2.2
1	B	152	VAL	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	734	LEU	2.2
1	A	155	ASN	2.2
1	B	622	ALA	2.2
1	A	28	GLU	2.1
1	B	613	GLY	2.1
1	B	628	PRO	2.1
1	A	368	ASN	2.1
1	B	535	SER	2.1
1	B	753	LEU	2.1
1	B	627	SER	2.1
1	A	228	ARG	2.1
1	A	647	ASN	2.1
1	B	691	GLU	2.1
1	A	734	LEU	2.0
1	A	343	ASP	2.0
1	B	154	PRO	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 monosaccharides 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, 95^{th} 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
2	GOL	A	801	6/6	0.94	0.18	29,32,36,36	0
2	GOL	B	801	6/6	0.94	0.16	34,40,43,47	0

6.5 Other polymers (i)

There are no such residues in this entry.