



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 23, 2024 – 11:22 PM EDT

PDB ID : 6Q9Z
Title : Crystal structure of the pathological G167R variant of calcium-free human gelsolin,
Authors : Boni, F.; Scalone, E.; Milani, M.; Eloise, M.; de Rosa, M.
Deposited on : 2018-12-18
Resolution : 3.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
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)	:	2.37.1

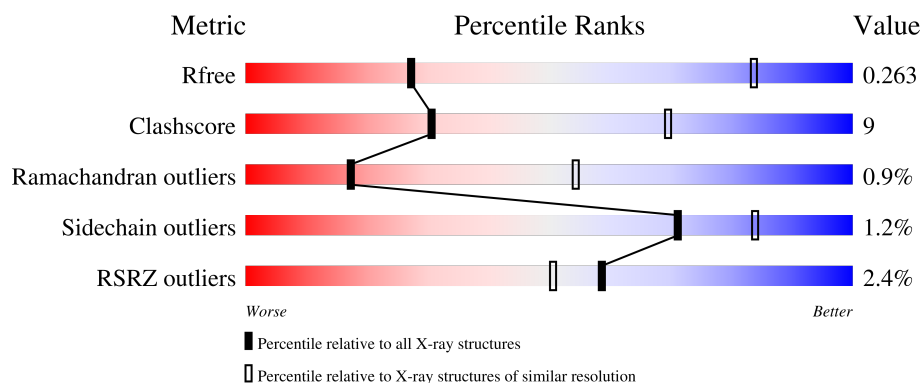
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.80 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	1212 (4.00-3.60)
Clashscore	141614	1288 (4.00-3.60)
Ramachandran outliers	138981	1243 (4.00-3.60)
Sidechain outliers	138945	1237 (4.00-3.60)
RSRZ outliers	127900	1121 (4.00-3.60)

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	778	
1	B	778	

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
3	GOL	B	802	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11169 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	708	Total	C	N	O	S	0	0	0
			5521	3484	962	1059	16			
1	B	716	Total	C	N	O	S	0	0	0
			5586	3526	980	1064	16			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-22	MET	-	initiating methionine	UNP P06396
A	-21	GLY	-	expression tag	UNP P06396
A	-20	SER	-	expression tag	UNP P06396
A	-19	SER	-	expression tag	UNP P06396
A	-18	HIS	-	expression tag	UNP P06396
A	-17	HIS	-	expression tag	UNP P06396
A	-16	HIS	-	expression tag	UNP P06396
A	-15	HIS	-	expression tag	UNP P06396
A	-14	HIS	-	expression tag	UNP P06396
A	-13	HIS	-	expression tag	UNP P06396
A	-12	SER	-	expression tag	UNP P06396
A	-11	SER	-	expression tag	UNP P06396
A	-10	GLY	-	expression tag	UNP P06396
A	-9	LEU	-	expression tag	UNP P06396
A	-8	VAL	-	expression tag	UNP P06396
A	-7	PRO	-	expression tag	UNP P06396
A	-6	ARG	-	expression tag	UNP P06396
A	-5	GLY	-	expression tag	UNP P06396
A	-4	SER	-	expression tag	UNP P06396
A	-3	HIS	-	expression tag	UNP P06396
A	-2	MET	-	expression tag	UNP P06396
A	-1	ALA	-	expression tag	UNP P06396
A	0	SER	-	expression tag	UNP P06396
A	167	ARG	GLY	engineered mutation	UNP P06396
B	-22	MET	-	initiating methionine	UNP P06396

Continued on next page...

Continued from previous page...

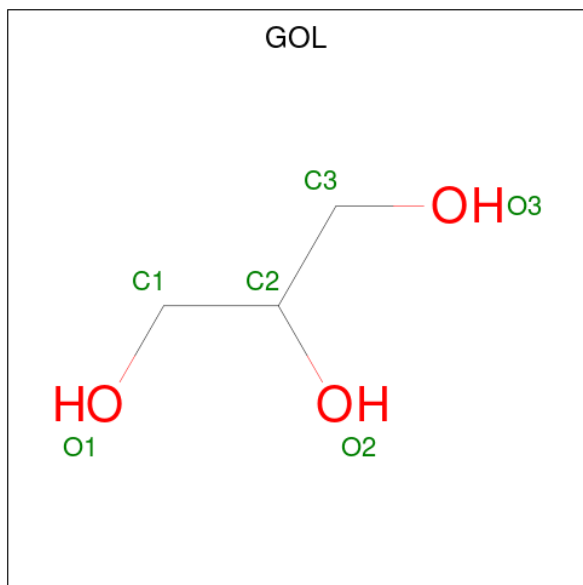
Chain	Residue	Modelled	Actual	Comment	Reference
B	-21	GLY	-	expression tag	UNP P06396
B	-20	SER	-	expression tag	UNP P06396
B	-19	SER	-	expression tag	UNP P06396
B	-18	HIS	-	expression tag	UNP P06396
B	-17	HIS	-	expression tag	UNP P06396
B	-16	HIS	-	expression tag	UNP P06396
B	-15	HIS	-	expression tag	UNP P06396
B	-14	HIS	-	expression tag	UNP P06396
B	-13	HIS	-	expression tag	UNP P06396
B	-12	SER	-	expression tag	UNP P06396
B	-11	SER	-	expression tag	UNP P06396
B	-10	GLY	-	expression tag	UNP P06396
B	-9	LEU	-	expression tag	UNP P06396
B	-8	VAL	-	expression tag	UNP P06396
B	-7	PRO	-	expression tag	UNP P06396
B	-6	ARG	-	expression tag	UNP P06396
B	-5	GLY	-	expression tag	UNP P06396
B	-4	SER	-	expression tag	UNP P06396
B	-3	HIS	-	expression tag	UNP P06396
B	-2	MET	-	expression tag	UNP P06396
B	-1	ALA	-	expression tag	UNP P06396
B	0	SER	-	expression tag	UNP P06396
B	167	ARG	GLY	engineered mutation	UNP P06396

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

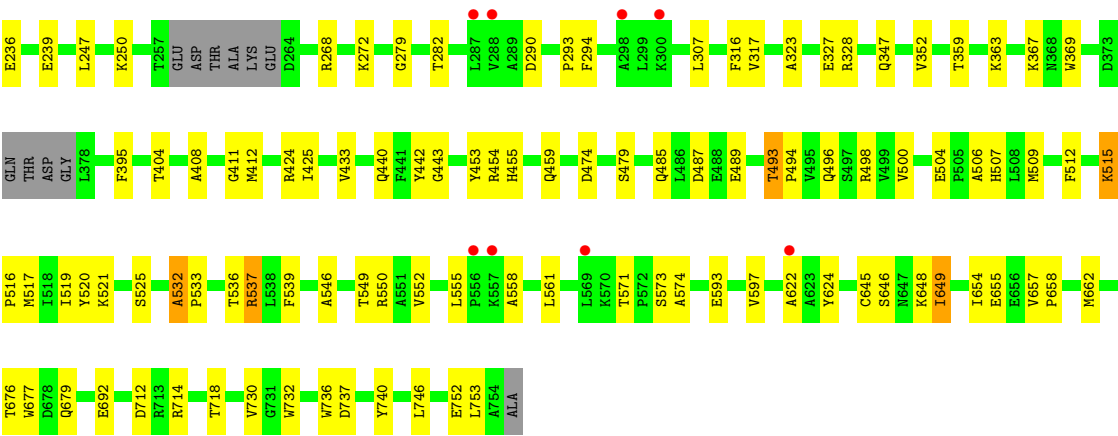
- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	19	Total O 19 19	0	0
4	B	21	Total O 21 21	0	0



4 Data and refinement statistics

Property	Value	Source
Space group	P 4 21 2	Depositor
Cell constants a, b, c, α , β , γ	170.91Å 170.91Å 151.62Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.47 – 3.80 48.47 – 3.80	Depositor EDS
% Data completeness (in resolution range)	99.9 (48.47-3.80) 99.9 (48.47-3.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 3.77Å)	Xtriage
Refinement program	PHENIX (1.10.1_2155: ???)	Depositor
R, R_{free}	0.216 , 0.264 0.220 , 0.263	Depositor DCC
R_{free} test set	1997 reflections (8.80%)	wwPDB-VP
Wilson B-factor (Å ²)	107.9	Xtriage
Anisotropy	0.296	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 85.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	11169	wwPDB-VP
Average B, all atoms (Å ²)	109.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% 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: GOL, SO4

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	0.26	0/5644	0.45	0/7643
1	B	0.27	1/5714 (0.0%)	0.45	0/7740
All	All	0.26	1/11358 (0.0%)	0.45	0/15383

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	555	LEU	C-N	6.92	1.47	1.34

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	5521	0	5355	119	0
1	B	5586	0	5448	86	0
2	A	5	0	0	0	0
2	B	5	0	0	0	0
3	B	12	0	16	0	0
4	A	19	0	0	0	0
4	B	21	0	0	0	0
All	All	11169	0	10819	201	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 9.

All (201) 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:595:LEU:HD23	1:A:602:PRO:HB3	1.64	0.80
1:A:268:ARG:HH22	1:A:707:ASP:HA	1.47	0.79
1:B:45:ARG:HH12	1:B:150:LYS:HA	1.51	0.76
1:A:168:ARG:O	1:A:170:VAL:N	2.21	0.73
1:A:33:LEU:H	1:A:33:LEU:HD13	1.53	0.73
1:B:532:ALA:HB1	1:B:533:PRO:HD2	1.71	0.72
1:B:323:ALA:O	1:B:328:ARG:NH1	2.24	0.70
1:A:268:ARG:HH11	1:A:654:ILE:HD11	1.57	0.69
1:A:389:ASN:ND2	1:A:660:GLU:OE2	2.26	0.69
1:A:302:GLU:HA	1:A:322:GLN:HB2	1.73	0.69
1:B:126:GLU:OE1	1:B:172:ARG:NH2	2.21	0.68
1:B:487:ASP:OD2	1:B:496:GLN:NE2	2.26	0.68
1:A:519:ILE:HD12	1:A:552:VAL:HG22	1.76	0.67
1:A:269:LYS:O	1:A:292:ASN:ND2	2.27	0.67
1:B:317:VAL:HB	1:B:352:VAL:HG12	1.76	0.66
1:B:519:ILE:HD12	1:B:552:VAL:HG22	1.77	0.65
1:A:509:MET:HE1	1:A:550:ARG:HH11	1.60	0.65
1:A:625:ARG:HE	1:A:626:THR:H	1.45	0.65
1:A:674:LEU:HB3	1:A:681:PHE:HB2	1.77	0.64
1:B:737:ASP:HB3	1:B:740:TYR:HB2	1.79	0.63
1:A:453:TYR:O	1:A:459:GLN:HB2	1.98	0.62
1:B:45:ARG:NH2	1:B:149:PHE:O	2.23	0.62
1:A:45:ARG:HH12	1:A:150:LYS:HA	1.66	0.60
1:B:167:ARG:NH2	1:B:655:GLU:OE2	2.34	0.60
1:B:294:PHE:HE2	1:B:307:LEU:HD13	1.67	0.60
1:B:500:VAL:HB	1:B:714:ARG:HB2	1.84	0.60
1:A:317:VAL:HG21	1:A:335:ALA:HB2	1.83	0.59
1:B:509:MET:HE1	1:B:550:ARG:HH11	1.67	0.59
1:A:750:MET:HE3	1:A:750:MET:HA	1.85	0.59
1:A:267:ASN:ND2	1:A:311:LYS:HB3	2.18	0.59
1:B:645:CYS:HB3	1:B:654:ILE:HG22	1.84	0.58
1:A:267:ASN:HD22	1:A:311:LYS:HB3	1.69	0.58
1:A:420:LYS:HE3	1:A:512:PHE:HD1	1.69	0.57
1:B:85:LEU:HD21	1:B:108:LEU:HD23	1.86	0.57
1:B:676:THR:HG22	1:B:677:TRP:H	1.70	0.57
1:A:200:TRP:CZ2	1:A:236:GLU:HG2	2.40	0.56
1:B:200:TRP:CZ2	1:B:236:GLU:HG2	2.39	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:504:GLU:HG2	1:B:517:MET:HE1	1.86	0.56
1:A:209:GLU:OE1	1:A:209:GLU:N	2.38	0.56
1:B:61:ASP:OD1	1:B:139:LYS:NZ	2.39	0.56
1:A:76:LEU:HB3	1:A:80:ASN:HB2	1.88	0.56
1:A:676:THR:HG22	1:A:677:TRP:H	1.71	0.56
1:B:573:SER:OG	1:B:574:ALA:N	2.38	0.56
1:B:45:ARG:NH1	1:B:150:LYS:HA	2.21	0.55
1:B:752:GLU:HG2	1:B:753:LEU:HD22	1.87	0.55
1:A:214:THR:HG22	1:A:218:LYS:HE3	1.89	0.55
1:B:498:ARG:NH1	1:B:500:VAL:HG23	2.21	0.55
1:A:516:PRO:HB3	1:A:549:THR:HG23	1.89	0.55
1:A:561:LEU:O	1:A:624:TYR:HB2	2.07	0.55
1:A:179:SER:HB2	1:B:177:PRO:HG3	1.89	0.54
1:A:177:PRO:HG3	1:B:179:SER:HB2	1.88	0.54
1:A:257:THR:HG22	1:A:258:GLU:H	1.72	0.54
1:B:118:GLN:HB2	1:B:352:VAL:HG22	1.88	0.54
1:B:57:ASN:O	1:B:223:ASN:HB3	2.08	0.54
1:A:221:ARG:NH2	1:A:753:LEU:HD23	2.22	0.54
1:A:500:VAL:HB	1:A:714:ARG:HB3	1.91	0.53
1:B:395:PHE:CE2	1:B:546:ALA:HA	2.43	0.53
1:A:517:MET:HE2	1:A:519:ILE:HD11	1.92	0.52
1:B:558:ALA:HB3	1:B:622:ALA:HB3	1.92	0.52
1:A:420:LYS:HE2	1:A:438:TYR:HD2	1.74	0.52
1:A:605:VAL:HG21	1:A:611:PRO:HD3	1.91	0.52
1:A:76:LEU:O	1:A:78:ASN:N	2.42	0.52
1:A:420:LYS:HE3	1:A:512:PHE:CD1	2.45	0.52
1:B:168:ARG:O	1:B:170:VAL:N	2.43	0.52
1:B:221:ARG:HG3	1:B:229:ALA:HB3	1.92	0.51
1:B:532:ALA:HB1	1:B:533:PRO:CD	2.40	0.51
1:A:487:ASP:OD2	1:A:496:GLN:NE2	2.43	0.51
1:A:371:ASP:OD1	1:A:371:ASP:N	2.44	0.51
1:B:424:ARG:HB2	1:B:433:VAL:HG22	1.92	0.51
1:A:625:ARG:HE	1:A:626:THR:N	2.08	0.51
1:B:506:ALA:HA	1:B:509:MET:HE3	1.93	0.51
1:B:679:GLN:NE2	1:B:718:THR:OG1	2.32	0.51
1:A:506:ALA:HA	1:A:509:MET:HE3	1.92	0.51
1:A:589:THR:O	1:A:592:GLN:HG2	2.11	0.51
1:A:634:LYS:NZ	1:A:678:ASP:OD2	2.44	0.50
1:A:395:PHE:CE2	1:A:546:ALA:HA	2.46	0.50
1:A:449:ILE:HB	1:A:464:TYR:HB2	1.94	0.50
1:B:537:ARG:HG3	1:B:539:PHE:HE1	1.76	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:LEU:HD21	1:A:108:LEU:HD23	1.93	0.50
1:A:424:ARG:HB2	1:A:433:VAL:HG22	1.94	0.49
1:A:425:ILE:HG13	1:A:479:SER:HB3	1.95	0.49
1:B:206:ASN:HB3	1:B:209:GLU:OE1	2.13	0.49
1:A:77:ARG:CG	1:A:378:LEU:HB3	2.42	0.49
1:A:221:ARG:HG3	1:A:229:ALA:HB3	1.93	0.49
1:B:307:LEU:HB3	1:B:316:PHE:HB2	1.94	0.49
1:B:363:LYS:HD2	1:B:369:TRP:CD1	2.48	0.49
1:A:168:ARG:HE	1:A:208:TYR:HD2	1.61	0.49
1:B:593:GLU:O	1:B:597:VAL:HG23	2.14	0.48
1:A:564:ASN:HD21	1:A:628:PRO:HA	1.79	0.48
1:A:509:MET:HE1	1:A:550:ARG:NH1	2.28	0.48
1:B:154:PRO:HD2	1:B:157:VAL:HG11	1.95	0.48
1:A:206:ASN:HB3	1:A:209:GLU:OE1	2.14	0.48
1:A:424:ARG:NH1	1:A:433:VAL:HG13	2.29	0.47
1:A:380:LEU:HD11	1:A:633:LYS:HB3	1.96	0.47
1:A:408:ALA:HB1	1:A:732:TRP:CE2	2.49	0.47
1:B:712:ASP:OD1	1:B:714:ARG:HG2	2.15	0.47
1:B:536:THR:HG22	1:B:571:THR:HA	1.97	0.47
1:B:112:LEU:O	1:B:115:ARG:HG2	2.15	0.47
1:A:591:ALA:O	1:A:595:LEU:HD13	2.16	0.46
1:B:207:ARG:HG3	1:B:746:LEU:HD22	1.96	0.46
1:A:426:GLU:HG2	1:A:526:ARG:HG2	1.97	0.46
1:B:272:LYS:HG2	1:B:290:ASP:HA	1.97	0.46
1:B:442:TYR:CE1	1:B:525:SER:HB3	2.50	0.46
1:A:293:PRO:HG3	1:A:710:ASN:HB2	1.98	0.46
1:A:347:GLN:OE1	1:A:347:GLN:N	2.47	0.46
1:A:537:ARG:HD2	1:A:553:GLU:OE1	2.16	0.46
1:B:135:LYS:HE3	1:B:662:MET:SD	2.55	0.46
1:A:239:GLU:HB2	1:A:244:LEU:HD21	1.97	0.46
1:B:183:PHE:HA	1:B:189:PHE:CZ	2.50	0.46
1:B:139:LYS:HE2	1:B:220:ILE:HG12	1.95	0.46
1:A:72:LYS:HE3	1:A:74:VAL:HG13	1.97	0.45
1:A:153:VAL:H	1:A:154:PRO:HD2	1.81	0.45
1:B:293:PRO:HB2	1:B:367:LYS:HE3	1.98	0.45
1:A:183:PHE:HA	1:A:189:PHE:CZ	2.51	0.45
1:A:558:ALA:HB3	1:A:622:ALA:HB3	1.99	0.45
1:A:201:CYS:O	1:A:235:GLU:HG3	2.16	0.45
1:A:749:ALA:HA	1:A:752:GLU:HB2	1.98	0.45
1:B:512:PHE:O	1:B:515:LYS:HB2	2.15	0.45
1:A:61:ASP:OD1	1:A:139:LYS:NZ	2.49	0.45

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:408:ALA:HB1	1:A:732:TRP:NE1	2.32	0.45
1:B:408:ALA:HB1	1:B:732:TRP:NE1	2.32	0.45
1:B:730:VAL:HA	1:B:736:TRP:CZ3	2.52	0.45
1:A:583:ALA:HA	1:A:587:GLU:OE2	2.17	0.45
1:B:347:GLN:OE1	1:B:347:GLN:N	2.48	0.44
1:B:180:TRP:NE1	1:B:247:LEU:HB3	2.31	0.44
1:B:359:THR:O	1:B:363:LYS:HG3	2.17	0.44
1:A:235:GLU:HG2	1:A:236:GLU:H	1.81	0.44
1:A:268:ARG:HB2	1:A:309:HIS:NE2	2.32	0.44
1:A:476:VAL:O	1:A:479:SER:OG	2.24	0.44
1:A:578:TRP:HE1	1:A:607:GLU:HA	1.82	0.44
1:B:154:PRO:HB2	1:B:157:VAL:HG13	1.99	0.44
1:B:648:LYS:HG3	1:B:649:ILE:HG13	2.00	0.44
1:A:411:GLY:O	1:A:507:HIS:HB2	2.16	0.44
1:A:549:THR:HG21	1:A:590:GLY:HA3	1.99	0.44
1:A:413:ASP:OD1	1:A:726:PRO:HB2	2.17	0.44
1:B:474:ASP:OD1	1:B:474:ASP:N	2.46	0.44
1:A:233:VAL:HG11	1:A:750:MET:CE	2.48	0.44
1:B:485:GLN:O	1:B:489:GLU:HG2	2.18	0.44
1:B:221:ARG:HH12	1:B:753:LEU:HB3	1.83	0.44
1:A:384:SER:OG	1:A:387:ILE:HG12	2.17	0.43
1:A:434:ASP:HB3	1:A:437:THR:OG1	2.17	0.43
1:B:347:GLN:H	1:B:347:GLN:CD	2.21	0.43
1:B:440:GLN:OE1	1:B:520:TYR:OH	2.26	0.43
1:A:33:LEU:H	1:A:33:LEU:CD1	2.27	0.43
1:A:77:ARG:CZ	1:A:392:ARG:HG3	2.48	0.43
1:B:218:LYS:HG2	1:B:222:ASP:OD2	2.19	0.43
1:A:169:ARG:HB3	1:A:169:ARG:NH2	2.33	0.43
1:A:474:ASP:OD1	1:A:474:ASP:N	2.52	0.43
1:A:179:SER:HB2	1:B:177:PRO:CG	2.49	0.43
1:A:536:THR:OG1	1:A:571:THR:HA	2.19	0.43
1:B:646:SER:HA	1:B:692:GLU:OE2	2.19	0.43
1:B:443:GLY:HA3	1:B:521:LYS:O	2.19	0.43
1:A:575:ALA:HB3	1:A:601:GLN:O	2.19	0.42
1:B:239:GLU:OE2	1:B:250:LYS:HD2	2.18	0.42
1:B:561:LEU:O	1:B:624:TYR:HB2	2.19	0.42
1:A:72:LYS:HB3	1:A:84:ASP:HB2	2.02	0.42
1:A:725:GLU:HB2	1:A:730:VAL:HG13	2.00	0.42
1:A:85:LEU:HA	1:A:85:LEU:HD23	1.74	0.42
1:B:425:ILE:HG13	1:B:479:SER:HB3	2.02	0.42
1:A:34:LYS:O	1:A:34:LYS:HD3	2.19	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:LEU:O	1:A:119:HIS:N	2.51	0.42
1:A:561:LEU:HD13	1:A:561:LEU:HA	1.90	0.42
1:A:77:ARG:HG2	1:A:378:LEU:HB3	2.02	0.42
1:A:499:VAL:HG21	1:A:504:GLU:HA	2.00	0.42
1:B:498:ARG:HH12	1:B:500:VAL:HG23	1.81	0.42
1:B:657:VAL:HA	1:B:658:PRO:HD3	1.90	0.42
1:B:37:LYS:HA	1:B:37:LYS:HD3	1.77	0.42
1:B:509:MET:HE1	1:B:550:ARG:NH1	2.34	0.42
1:A:45:ARG:HB2	1:A:68:TYR:CE2	2.55	0.41
1:A:198:HIS:HA	1:A:232:HIS:O	2.20	0.41
1:A:268:ARG:NH2	1:A:707:ASP:HA	2.25	0.41
1:A:412:MET:SD	1:A:507:HIS:HA	2.60	0.41
1:A:587:GLU:H	1:A:587:GLU:HG2	1.70	0.41
1:B:209:GLU:OE1	1:B:209:GLU:N	2.48	0.41
1:A:64:THR:HB	1:A:92:GLU:HG2	2.02	0.41
1:A:423:TRP:CD1	1:A:430:LYS:HB2	2.55	0.41
1:A:425:ILE:HG13	1:A:479:SER:CB	2.50	0.41
1:A:675:ASP:OD1	1:A:680:VAL:HG22	2.19	0.41
1:B:279:GLY:HA3	1:B:327:GLU:OE1	2.21	0.41
1:B:516:PRO:HB3	1:B:549:THR:OG1	2.19	0.41
1:B:736:TRP:HD1	1:B:737:ASP:N	2.17	0.41
1:A:314:LYS:HE2	1:A:349:GLN:OE1	2.20	0.41
1:A:666:LEU:HD22	1:A:672:MET:HG3	2.02	0.41
1:B:411:GLY:O	1:B:507:HIS:HB2	2.20	0.41
1:A:392:ARG:HH21	1:A:636:ASP:CG	2.24	0.41
1:B:408:ALA:HB1	1:B:732:TRP:CD1	2.55	0.41
1:A:246:VAL:O	1:B:142:LYS:HD2	2.20	0.41
1:A:292:ASN:HB3	1:A:293:PRO:HD3	2.02	0.41
1:A:512:PHE:O	1:A:515:LYS:HB2	2.21	0.41
1:A:612:ASP:OD2	1:A:612:ASP:N	2.54	0.41
1:B:412:MET:SD	1:B:507:HIS:HA	2.60	0.41
1:A:566:ALA:HB3	1:A:587:GLU:HB2	2.02	0.40
1:B:493:THR:H	1:B:494:PRO:CD	2.35	0.40
1:A:267:ASN:ND2	1:A:267:ASN:O	2.54	0.40
1:A:565:ASP:HA	1:A:587:GLU:OE1	2.21	0.40
1:B:453:TYR:O	1:B:459:GLN:HB2	2.21	0.40
1:A:568:VAL:HG21	1:A:594:LEU:HD23	2.03	0.40
1:B:404:THR:HG22	1:B:455:HIS:HE1	1.86	0.40
1:A:266:ALA:HB1	1:A:706:THR:HG21	2.03	0.40
1:A:380:LEU:CD1	1:A:633:LYS:HB3	2.51	0.40
1:A:383:LEU:HD21	1:A:388:ALA:HA	2.03	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	694/778 (89%)	643 (93%)	44 (6%)	7 (1%)	15	52
1	B	710/778 (91%)	661 (93%)	44 (6%)	5 (1%)	22	60
All	All	1404/1556 (90%)	1304 (93%)	88 (6%)	12 (1%)	17	54

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	169	ARG
1	A	30	PRO
1	A	77	ARG
1	A	372	PRO
1	B	169	ARG
1	B	532	ALA
1	B	649	ILE
1	A	79	GLY
1	B	282	THR
1	A	493	THR
1	B	493	THR
1	A	153	VAL

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	578/630 (92%)	571 (99%)	7 (1%)	71	84
1	B	584/630 (93%)	577 (99%)	7 (1%)	71	84
All	All	1162/1260 (92%)	1148 (99%)	14 (1%)	71	84

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	33	LEU
1	A	169	ARG
1	A	268	ARG
1	A	498	ARG
1	A	593	GLU
1	A	612	ASP
1	A	641	ARG
1	B	76	LEU
1	B	169	ARG
1	B	218	LYS
1	B	268	ARG
1	B	454	ARG
1	B	515	LYS
1	B	537	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	267	ASN
1	B	467	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 [i](#)

4 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	$\# Z > 2$
3	GOL	B	801	-	5,5,5	0.36	0	5,5,5	0.34	0
3	GOL	B	802	-	5,5,5	0.37	0	5,5,5	0.34	0
2	SO4	A	801	-	4,4,4	0.23	0	6,6,6	0.16	0
2	SO4	B	803	-	4,4,4	0.24	0	6,6,6	0.06	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
3	GOL	B	801	-	-	0/4/4/4	-
3	GOL	B	802	-	-	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.

No monomer is involved in short contacts.

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	708/778 (91%)	0.10	26 (3%) 41 34	82, 105, 173, 256	0
1	B	716/778 (92%)	-0.04	8 (1%) 80 74	80, 100, 157, 228	0
All	All	1424/1556 (91%)	0.03	34 (2%) 59 50	80, 103, 167, 256	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	576	TYR	4.6
1	A	575	ALA	4.0
1	A	568	VAL	3.8
1	A	298	ALA	3.6
1	A	574	ALA	3.4
1	A	622	ALA	3.3
1	A	303	ASP	3.0
1	A	558	ALA	3.0
1	B	622	ALA	2.9
1	A	754	ALA	2.9
1	A	300	LYS	2.7
1	A	557	LYS	2.6
1	B	556	PRO	2.6
1	A	571	THR	2.5
1	B	557	LYS	2.5
1	A	304	CYS	2.4
1	A	273	LEU	2.4
1	B	569	LEU	2.4
1	B	287	LEU	2.4
1	A	154	PRO	2.3
1	A	305	PHE	2.3
1	A	274	TYR	2.3
1	A	151	HIS	2.3
1	A	612	ASP	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	556	PRO	2.3
1	A	299	LEU	2.3
1	B	288	VAL	2.2
1	A	573	SER	2.2
1	B	300	LYS	2.2
1	A	539	PHE	2.2
1	B	298	ALA	2.2
1	A	752	GLU	2.2
1	A	276	VAL	2.2
1	A	611	PRO	2.1

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, 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	GOL	B	802	6/6	0.56	0.78	92,106,107,107	0
2	SO4	A	801	5/5	0.78	0.30	133,134,142,158	0
3	GOL	B	801	6/6	0.91	0.55	64,64,64,68	0
2	SO4	B	803	5/5	0.93	0.63	145,146,146,147	0

6.5 Other polymers [i](#)

There are no such residues in this entry.