



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 01:16 pm BST

PDB ID : 2E0Y
Title : Crystal structure of the samarium derivative of mature gamma-glutamyltranspeptidase from Escherichia coli
Authors : Okada, T.; Wada, K.; Fukuyama, K.
Deposited on : 2006-10-16
Resolution : 2.02 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

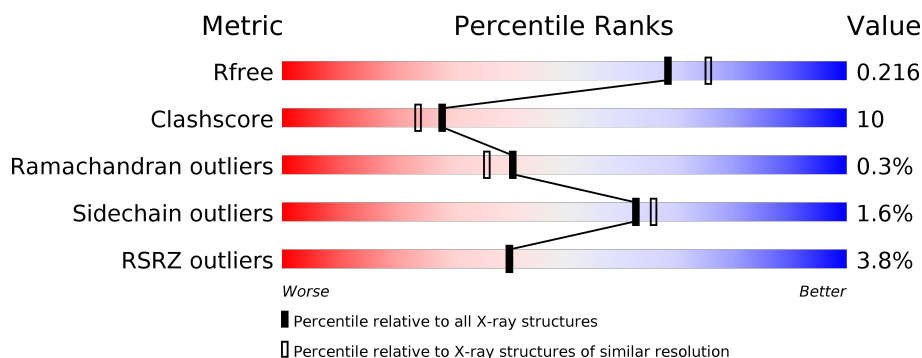
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.02 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	10434 (2.04-2.00)
Clashscore	141614	11643 (2.04-2.00)
Ramachandran outliers	138981	11493 (2.04-2.00)
Sidechain outliers	138945	11492 (2.04-2.00)
RSRZ outliers	127900	10220 (2.04-2.00)

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 on 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	366	<div> <div>4%</div> <div> <div></div> <div>80%</div> <div>17%</div> <div>• •</div> </div> </div>
1	C	366	<div> <div>4%</div> <div> <div></div> <div>82%</div> <div>15%</div> <div>• •</div> </div> </div>
2	B	190	<div> <div>3%</div> <div> <div></div> <div>78%</div> <div>16%</div> <div>• 5%</div> </div> </div>
2	D	190	<div> <div>3%</div> <div> <div></div> <div>76%</div> <div>18%</div> <div>• 5%</div> </div> </div>

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 8706 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.

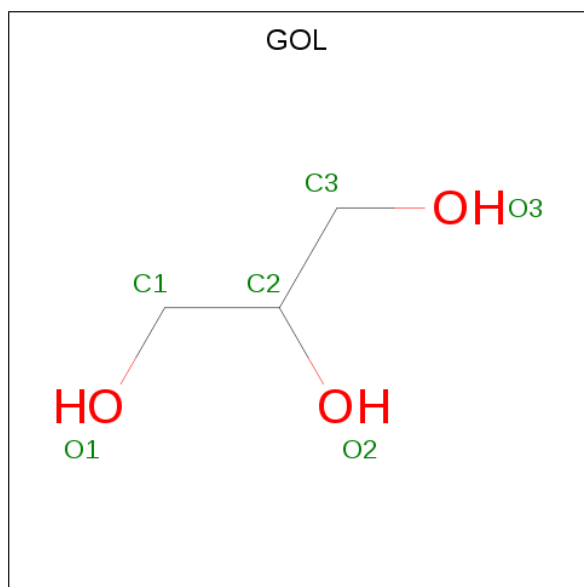
- Molecule 1 is a protein called Gamma-glutamyltranspeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	359	Total	C	N	O	Se	0	0	0
			2716	1717	456	532	11			
1	C	358	Total	C	N	O	Se	0	0	0
			2711	1714	455	531	11			

- Molecule 2 is a protein called Gamma-glutamyltranspeptidase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	180	Total	C	N	O	Se	0	0	0
			1336	834	227	270	5			
2	D	180	Total	C	N	O	Se	0	0	0
			1336	834	227	270	5			

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



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

- Molecule 4 is SAMARIUM (III) ION (three-letter code: SM) (formula: Sm).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Sm 1 1	0	0
4	D	1	Total Sm 1 1	0	0

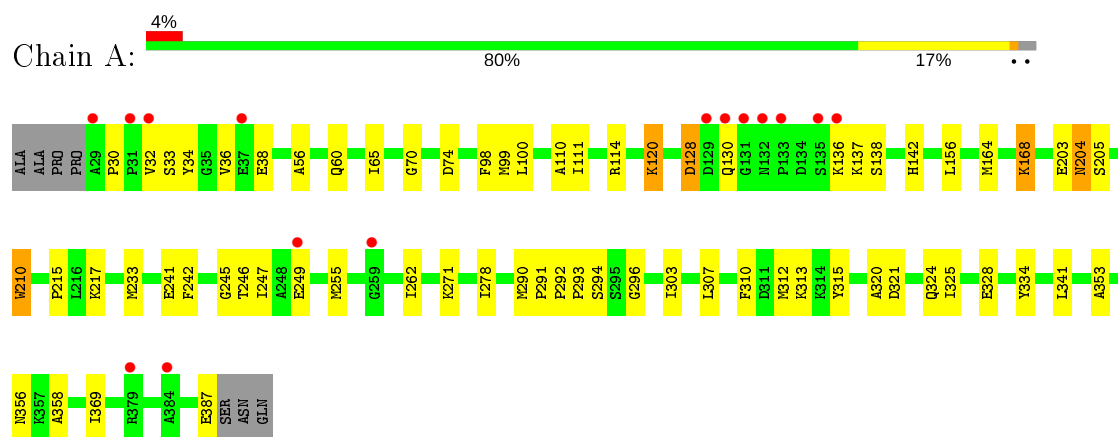
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	187	Total O 187 187	0	0
5	B	91	Total O 91 91	0	0
5	C	209	Total O 209 209	0	0
5	D	88	Total O 88 88	0	0

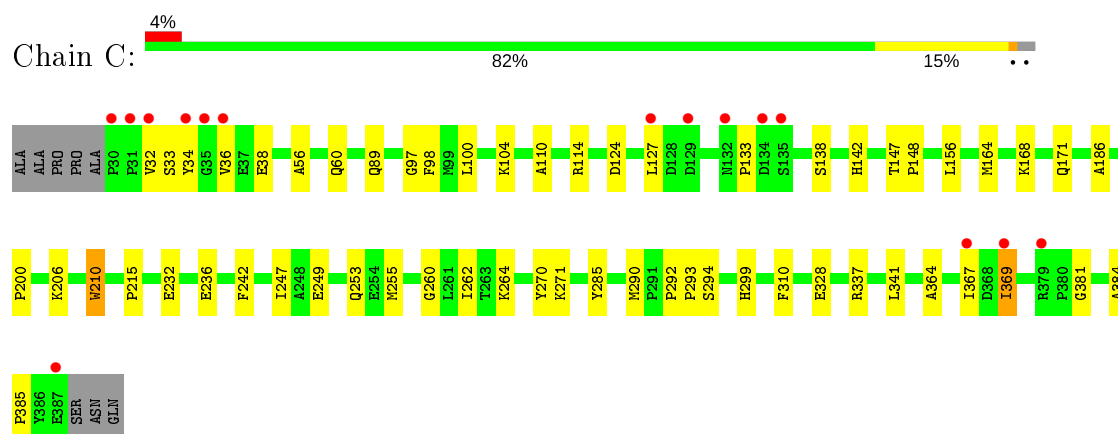
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

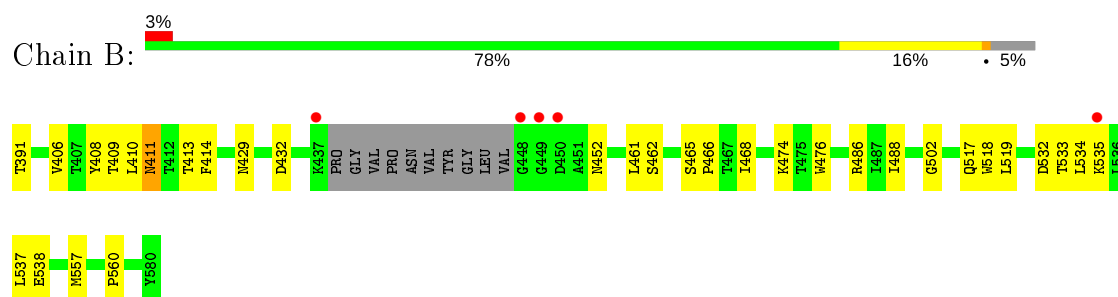
• Molecule 1: Gamma-glutamyltranspeptidase



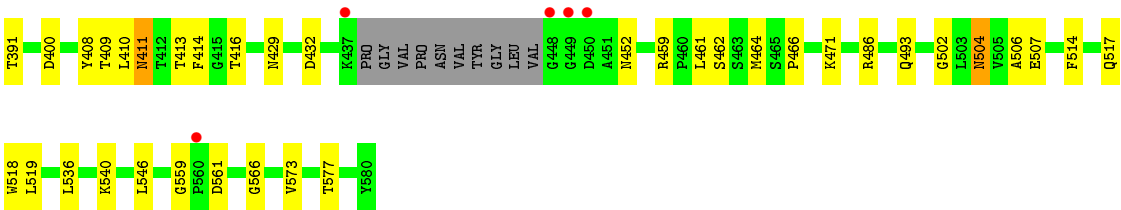
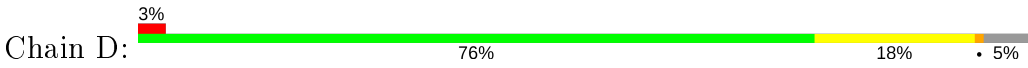
• Molecule 1: Gamma-glutamyltranspeptidase



• Molecule 2: Gamma-glutamyltranspeptidase



● Molecule 2: Gamma-glutamyltranspeptidase



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	78.92Å 125.90Å 128.90Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.02 46.40 – 2.02	Depositor EDS
% Data completeness (in resolution range)	98.8 (50.00-2.02) 99.0 (46.40-2.02)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.22 (at 2.01Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.195 , 0.224 0.189 , 0.216	Depositor DCC
R_{free} test set	4282 reflections (5.06%)	wwPDB-VP
Wilson B-factor (Å ²)	25.1	Xtriage
Anisotropy	0.345	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 44.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.011 for -h,l,k	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8706	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.66% 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: GOL, SM

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	0.30	0/2761	0.57	0/3721
1	C	0.30	0/2756	0.57	0/3713
2	B	0.30	0/1354	0.66	0/1830
2	D	0.30	0/1354	0.67	0/1830
All	All	0.30	0/8225	0.60	0/11094

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

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	2716	0	2680	74	0
1	C	2711	0	2676	48	0
2	B	1336	0	1318	30	0
2	D	1336	0	1318	35	0
3	A	12	0	16	0	0
3	B	6	0	8	0	0
3	C	6	0	8	1	0
3	D	6	0	8	0	0
4	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	1	0	0	0	0
5	A	187	0	0	1	0
5	B	91	0	0	1	0
5	C	209	0	0	2	0
5	D	88	0	0	0	0
All	All	8706	0	8032	158	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 (158) 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:74:ASP:HB3	1:A:164:MSE:HE3	1.22	1.18
1:A:70:GLY:HA3	1:A:164:MSE:HE2	1.25	1.15
1:A:65:ILE:HG23	1:A:164:MSE:HE1	1.27	1.12
1:A:74:ASP:CB	1:A:164:MSE:HE3	1.90	1.01
1:A:70:GLY:CA	1:A:164:MSE:HE2	1.94	0.96
1:C:290:MSE:HE3	2:D:466:PRO:HG2	1.50	0.93
1:A:70:GLY:HA3	1:A:164:MSE:CE	2.00	0.92
1:A:65:ILE:CG2	1:A:164:MSE:HE1	2.01	0.91
1:A:312:MSE:HA	1:A:312:MSE:HE2	1.55	0.88
1:C:369:ILE:HD13	1:C:369:ILE:H	1.41	0.85
1:A:310:PHE:HB3	1:A:312:MSE:HE3	1.62	0.81
2:D:493:GLN:HE22	2:D:514:PHE:H	1.25	0.80
1:A:120:LYS:HZ3	1:A:120:LYS:HB2	1.50	0.77
1:C:164:MSE:HE2	1:C:168:LYS:HB3	1.67	0.74
1:A:30:PRO:HD3	1:A:313:LYS:HE2	1.69	0.74
1:A:310:PHE:CB	1:A:312:MSE:HE3	2.20	0.72
1:A:36:VAL:HB	1:A:38:GLU:OE2	1.90	0.71
1:C:249:GLU:OE1	1:C:264:LYS:HD2	1.90	0.70
1:A:56:ALA:O	1:A:60:GLN:HG3	1.92	0.69
1:A:98:PHE:CD1	1:A:290:MSE:HE2	2.28	0.69
2:D:504:ASN:HD22	2:D:506:ALA:H	1.44	0.65
1:A:204:ASN:HD22	1:A:205:SER:H	1.46	0.64
1:A:296:GLY:HA2	2:B:488:ILE:HG12	1.80	0.62
2:B:411:ASN:HB3	2:B:429:ASN:OD1	1.99	0.62
1:A:32:VAL:HG12	1:A:33:SER:N	2.14	0.62
1:A:353:ALA:HB2	1:A:387:GLU:HG2	1.83	0.61
2:B:532:ASP:O	2:B:535:LYS:HG2	2.01	0.61
2:D:411:ASN:HB3	2:D:429:ASN:OD1	1.99	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:504:ASN:ND2	2:D:507:GLU:H	1.98	0.60
1:A:136:LYS:HD3	1:A:136:LYS:C	2.20	0.60
1:C:369:ILE:CD1	1:C:369:ILE:H	2.14	0.60
2:B:391:THR:HA	2:B:409:THR:HB	1.86	0.58
1:A:321:ASP:CG	1:A:369:ILE:HD11	2.25	0.57
1:C:255:MSE:HG3	1:C:262:ILE:HB	1.87	0.56
2:D:504:ASN:ND2	2:D:506:ALA:H	2.03	0.56
1:C:290:MSE:HE3	2:D:466:PRO:CG	2.28	0.56
1:A:120:LYS:H	1:A:120:LYS:NZ	2.02	0.56
2:D:504:ASN:C	2:D:504:ASN:HD22	2.09	0.55
1:A:138:SER:O	2:B:432:ASP:HB2	2.06	0.55
2:B:533:THR:O	2:B:537:LEU:HD23	2.06	0.55
1:C:369:ILE:HD13	1:C:369:ILE:N	2.15	0.55
1:A:30:PRO:CD	1:A:313:LYS:HE2	2.37	0.54
1:A:38:GLU:CD	1:A:38:GLU:H	2.10	0.54
2:D:504:ASN:HD22	2:D:506:ALA:N	2.04	0.54
1:C:142:HIS:HD2	1:C:255:MSE:HE2	1.72	0.54
1:C:293:PRO:O	2:D:461:LEU:HD12	2.08	0.54
1:C:328:GLU:HB3	1:C:367:ILE:HG22	1.90	0.54
1:A:278:ILE:HG12	1:A:291:PRO:HB3	1.88	0.53
1:A:255:MSE:HG3	1:A:262:ILE:HB	1.90	0.53
2:B:518:TRP:O	2:B:519:LEU:HD12	2.09	0.53
1:A:204:ASN:HD22	1:A:205:SER:N	2.06	0.53
1:A:296:GLY:HA3	2:B:465:SER:OG	2.10	0.52
2:B:518:TRP:CD2	2:B:519:LEU:HD13	2.45	0.52
1:C:341:LEU:HB3	2:D:461:LEU:HD21	1.92	0.51
1:A:312:MSE:HE2	1:A:315:TYR:HD2	1.76	0.51
1:A:32:VAL:CG1	1:A:33:SER:N	2.74	0.51
1:A:204:ASN:N	1:A:204:ASN:HD22	2.09	0.50
1:C:56:ALA:O	1:C:60:GLN:HG3	2.12	0.50
1:C:200:PRO:O	1:C:206:LYS:HE3	2.11	0.50
2:D:518:TRP:CD2	2:D:519:LEU:HD13	2.46	0.50
1:C:364:ALA:O	1:C:367:ILE:HG12	2.11	0.50
1:C:232:GLU:O	1:C:236:GLU:HG3	2.12	0.49
1:C:384:ALA:N	1:C:385:PRO:HD2	2.27	0.49
2:D:517:GLN:O	2:D:518:TRP:HB3	2.12	0.49
2:D:413:THR:O	2:D:414:PHE:HB2	2.12	0.49
1:A:241:GLU:OE2	1:A:246:THR:HB	2.13	0.49
1:A:136:LYS:HD3	1:A:137:LYS:N	2.27	0.49
1:C:89:GLN:HB2	2:D:413:THR:HG23	1.93	0.49
1:A:98:PHE:CG	1:A:290:MSE:HE2	2.48	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:312:MSE:HA	1:A:312:MSE:CE	2.36	0.49
1:A:111:ILE:HD12	1:A:156:LEU:HD22	1.95	0.49
1:C:97:GLY:HA2	2:D:464:MSE:HE1	1.94	0.49
1:A:99:MSE:HG3	2:B:406:VAL:HG22	1.95	0.48
2:B:452:ASN:HD21	2:B:461:LEU:H	1.60	0.48
1:C:186:ALA:HB2	2:D:573:VAL:HG23	1.95	0.48
1:A:312:MSE:HE1	1:A:325:ILE:HD12	1.96	0.48
1:C:98:PHE:HB3	1:C:290:MSE:SE	2.64	0.48
1:A:156:LEU:C	1:A:156:LEU:HD23	2.34	0.48
2:B:413:THR:O	2:B:414:PHE:HB2	2.14	0.47
1:A:128:ASP:OD1	1:A:130:GLN:N	2.48	0.47
1:A:320:ALA:HA	2:B:537:LEU:HD21	1.95	0.47
1:C:32:VAL:HG12	1:C:33:SER:N	2.29	0.47
1:C:33:SER:HB3	2:D:502:GLY:O	2.14	0.47
1:A:100:LEU:HD23	2:B:468:ILE:HG13	1.95	0.47
1:A:290:MSE:HG2	2:B:466:PRO:HG2	1.97	0.47
1:A:34:TYR:HA	2:B:557:MSE:CE	2.44	0.47
1:C:381:GLY:HA2	3:C:704:GOL:H11	1.97	0.47
1:A:245:GLY:O	1:A:249:GLU:HG3	2.15	0.46
1:A:210:TRP:CH2	1:A:215:PRO:HB3	2.50	0.46
1:A:120:LYS:H	1:A:120:LYS:HZ2	1.63	0.46
1:A:242:PHE:HA	1:A:247:ILE:HB	1.96	0.46
1:C:114:ARG:CZ	2:D:462:SER:HB2	2.46	0.46
1:A:303:ILE:O	1:A:307:LEU:HG	2.16	0.46
2:B:517:GLN:O	2:B:518:TRP:HB3	2.14	0.46
1:C:292:PRO:HA	1:C:294:SER:N	2.31	0.46
1:C:310:PHE:CE2	1:C:367:ILE:HD11	2.51	0.46
1:A:32:VAL:HG13	5:B:1108:HOH:O	2.16	0.46
1:A:293:PRO:O	2:B:461:LEU:HD12	2.16	0.46
1:A:356:ASN:OD1	1:A:358:ALA:N	2.47	0.45
1:C:310:PHE:CZ	1:C:367:ILE:HD11	2.51	0.45
5:C:1070:HOH:O	2:D:459:ARG:HD2	2.15	0.45
2:D:546:LEU:HD13	2:D:546:LEU:C	2.36	0.45
1:A:98:PHE:HB3	1:A:290:MSE:SE	2.65	0.45
1:A:334:TYR:CD2	2:B:517:GLN:HA	2.52	0.45
2:B:410:LEU:O	2:B:411:ASN:HB3	2.17	0.45
1:C:34:TYR:CE2	2:D:559:GLY:HA2	2.52	0.45
1:C:255:MSE:SE	1:C:260:GLY:HA3	2.67	0.45
2:D:452:ASN:HD21	2:D:461:LEU:H	1.65	0.45
1:C:156:LEU:C	1:C:156:LEU:HD23	2.37	0.44
1:C:110:ALA:HB1	1:C:290:MSE:SE	2.68	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:504:ASN:ND2	2:D:506:ALA:N	2.63	0.44
2:B:534:LEU:O	2:B:538:GLU:HG3	2.17	0.44
1:C:210:TRP:CH2	1:C:215:PRO:HB3	2.53	0.44
1:A:114:ARG:CZ	2:B:462:SER:HB2	2.48	0.44
1:A:168:LYS:NZ	1:A:168:LYS:HB2	2.33	0.44
1:A:312:MSE:CA	1:A:312:MSE:HE2	2.36	0.44
2:D:504:ASN:HD21	2:D:506:ALA:HB3	1.83	0.44
1:C:328:GLU:HB3	1:C:367:ILE:CG2	2.47	0.43
1:C:104:LYS:HD2	2:D:400:ASP:HB3	2.00	0.43
1:C:270:TYR:CD2	1:C:271:LYS:N	2.87	0.43
2:D:504:ASN:ND2	2:D:504:ASN:C	2.71	0.43
2:D:410:LEU:O	2:D:411:ASN:HB3	2.18	0.43
1:A:292:PRO:HA	1:A:294:SER:N	2.33	0.43
1:A:33:SER:HB3	2:B:502:GLY:O	2.18	0.43
1:C:32:VAL:CG1	1:C:33:SER:N	2.82	0.43
1:C:138:SER:O	2:D:432:ASP:HB2	2.19	0.43
2:B:474:LYS:HZ1	2:B:560:PRO:CB	2.32	0.43
1:C:299:HIS:NE2	1:C:337:ARG:CB	2.82	0.43
1:A:30:PRO:HD3	1:A:313:LYS:CE	2.44	0.42
1:A:341:LEU:HB3	2:B:461:LEU:HD21	2.01	0.42
1:C:36:VAL:HG23	1:C:38:GLU:HG2	2.01	0.42
2:D:566:GLY:HA3	2:D:577:THR:HG21	2.01	0.42
1:C:124:ASP:HB3	1:C:127:LEU:HD12	2.02	0.42
1:A:204:ASN:ND2	1:A:205:SER:H	2.16	0.42
2:B:452:ASN:ND2	2:B:461:LEU:H	2.17	0.42
1:C:264:LYS:HB3	1:C:264:LYS:HE2	1.90	0.42
1:A:110:ALA:HB1	1:A:290:MSE:SE	2.69	0.42
1:A:204:ASN:ND2	1:A:205:SER:N	2.68	0.42
1:C:127:LEU:HD23	1:C:133:PRO:HA	2.02	0.41
1:A:34:TYR:HA	2:B:557:MSE:HE1	2.01	0.41
1:C:242:PHE:HA	1:C:247:ILE:HB	2.02	0.41
1:A:156:LEU:HD23	1:A:156:LEU:O	2.21	0.41
1:A:290:MSE:CG	2:B:466:PRO:HD2	2.51	0.41
1:C:100:LEU:HD12	1:C:100:LEU:HA	1.94	0.41
2:D:536:LEU:O	2:D:540:LYS:HG3	2.21	0.41
1:A:215:PRO:HG2	1:A:217:LYS:HZ3	1.85	0.41
2:B:410:LEU:O	2:B:411:ASN:CB	2.68	0.41
1:C:285:TYR:HE1	2:D:471:LYS:HD2	1.86	0.41
1:A:142:HIS:HD2	1:A:255:MSE:HE2	1.86	0.40
1:C:171:GLN:NE2	5:C:1123:HOH:O	2.52	0.40
2:D:416:THR:HG22	2:D:429:ASN:HB3	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:32:VAL:HG11	2:B:476:TRP:HZ2	1.87	0.40
2:D:493:GLN:HE22	2:D:514:PHE:N	2.04	0.40
1:A:203:GLU:HG2	5:A:1170:HOH:O	2.21	0.40
1:A:233:MSE:HB3	1:A:241:GLU:HG2	2.03	0.40
1:A:324:GLN:O	1:A:328:GLU:HG2	2.21	0.40
1:C:147:THR:HA	1:C:148:PRO:HD3	1.95	0.40
2:D:391:THR:HA	2:D:409:THR:HB	2.02	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	357/366 (98%)	346 (97%)	10 (3%)	1 (0%)	41	36
1	C	356/366 (97%)	349 (98%)	7 (2%)	0	100	100
2	B	176/190 (93%)	170 (97%)	5 (3%)	1 (1%)	25	18
2	D	176/190 (93%)	171 (97%)	4 (2%)	1 (1%)	25	18
All	All	1065/1112 (96%)	1036 (97%)	26 (2%)	3 (0%)	41	36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	411	ASN
2	D	411	ASN
1	A	128	ASP

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	282/276 (102%)	277 (98%)	5 (2%)	59	61
1	C	282/276 (102%)	279 (99%)	3 (1%)	73	77
2	B	146/149 (98%)	144 (99%)	2 (1%)	67	70
2	D	146/149 (98%)	142 (97%)	4 (3%)	44	44
All	All	856/850 (101%)	842 (98%)	14 (2%)	62	66

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

Mol	Chain	Res	Type
1	A	120	LYS
1	A	168	LYS
1	A	204	ASN
1	A	210	TRP
1	A	271	LYS
2	B	408	TYR
2	B	486	ARG
1	C	210	TRP
1	C	253	GLN
1	C	369	ILE
2	D	408	TYR
2	D	486	ARG
2	D	504	ASN
2	D	561	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	ASN
1	A	201	ASN
1	A	204	ASN
1	A	250	GLN
1	A	253	GLN
2	B	452	ASN
1	C	107	ASN
1	C	132	ASN
1	C	171	GLN

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Mol	Chain	Res	Type
1	C	201	ASN
1	C	237	ASN
1	C	256	GLN
2	D	452	ASN
2	D	493	GLN
2	D	497	ASN
2	D	504	ASN

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 7 ligands modelled in this entry, 2 are monoatomic - leaving 5 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	$\# Z > 2$
3	GOL	B	702	-	5,5,5	0.37	0	5,5,5	0.22	0
3	GOL	A	701	-	5,5,5	0.39	0	5,5,5	0.24	0
3	GOL	A	705	-	5,5,5	0.27	0	5,5,5	0.26	0
3	GOL	D	703	-	5,5,5	0.34	0	5,5,5	0.26	0
3	GOL	C	704	-	5,5,5	0.31	0	5,5,5	0.26	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	702	-	-	0/4/4/4	-
3	GOL	A	701	-	-	0/4/4/4	-
3	GOL	A	705	-	-	0/4/4/4	-
3	GOL	D	703	-	-	0/4/4/4	-
3	GOL	C	704	-	-	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	704	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	348/366 (95%)	0.07	15 (4%) 35 34	16, 26, 47, 70	0
1	C	347/366 (94%)	0.16	15 (4%) 35 34	14, 26, 48, 71	0
2	B	175/190 (92%)	0.13	5 (2%) 51 51	16, 23, 37, 68	0
2	D	175/190 (92%)	0.15	5 (2%) 51 51	15, 22, 41, 67	0
All	All	1045/1112 (93%)	0.12	40 (3%) 40 40	14, 25, 47, 71	0

All (40) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	448	GLY	8.9
2	D	448	GLY	8.8
1	C	32	VAL	5.9
1	C	35	GLY	5.6
2	D	560	PRO	5.5
1	C	31	PRO	4.8
2	D	449	GLY	4.8
1	A	129	ASP	4.7
1	A	131	GLY	4.5
1	C	36	VAL	4.1
1	A	130	GLN	4.0
1	C	30	PRO	3.9
2	D	450	ASP	3.8
2	B	449	GLY	3.8
1	C	387	GLU	3.7
1	A	132	ASN	3.7
1	C	34	TYR	3.7
1	A	136	LYS	3.6
2	B	437	LYS	3.6
1	A	133	PRO	3.6
1	A	32	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
1	A	29	ALA	3.0
1	C	132	ASN	3.0
1	C	134	ASP	2.8
1	A	259	GLY	2.7
2	B	450	ASP	2.6
1	A	249	GLU	2.6
2	D	437	LYS	2.5
1	C	129	ASP	2.5
1	C	367	ILE	2.4
1	C	379	ARG	2.4
1	C	135	SER	2.4
1	A	384	ALA	2.3
1	C	369	ILE	2.3
2	B	535	LYS	2.2
1	A	135	SER	2.2
1	A	379	ARG	2.1
1	A	37	GLU	2.0
1	A	31	PRO	2.0
1	C	127	LEU	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 carbohydrates 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(\AA^2)	Q<0.9
3	GOL	A	701	6/6	0.76	0.21	46,47,48,48	0
3	GOL	A	705	6/6	0.88	0.22	53,54,55,56	0
3	GOL	C	704	6/6	0.89	0.25	44,46,46,46	0
3	GOL	D	703	6/6	0.94	0.15	38,40,42,43	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
3	GOL	B	702	6/6	0.94	0.14	30,37,38,39	0
4	SM	B	601	1/1	1.00	0.03	38,38,38,38	0
4	SM	D	602	1/1	1.00	0.02	35,35,35,35	0

6.5 Other polymers [i](#)

There are no such residues in this entry.