



Full wwPDB X-ray Structure Validation Report ⓘ

Dec 15, 2024 – 09:45 PM EST

PDB ID : 2O99
Title : The crystal structure of E.coli IclR C-terminal fragment in complex with glyoxylate
Authors : Lunin, V.V.; Ezersky, A.; Evdokimova, E.; Kudritska, M.; Savchenko, A.
Deposited on : 2006-12-13
Resolution : 1.70 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)	:	1.21
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.40

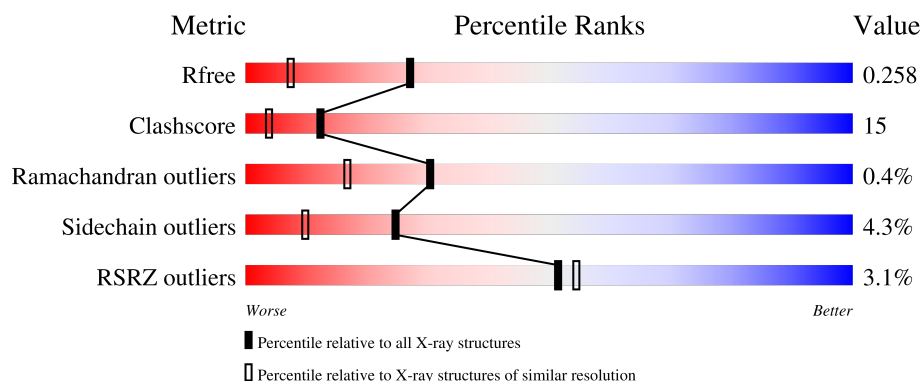
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 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}	164625	5161 (1.70-1.70)
Clashscore	180529	5671 (1.70-1.70)
Ramachandran outliers	177936	5594 (1.70-1.70)
Sidechain outliers	177891	5594 (1.70-1.70)
RSRZ outliers	164620	5159 (1.70-1.70)

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	182	<div> <div>4%</div> <div> <div></div> <div>79%</div> <div>16%</div> <div>..</div> </div> </div>
1	B	182	<div> <div>3%</div> <div> <div></div> <div>74%</div> <div>22%</div> <div>..</div> </div> </div>
1	C	182	<div> <div>4%</div> <div> <div></div> <div>79%</div> <div>18%</div> <div>..</div> </div> </div>
1	D	182	<div> <div>%</div> <div> <div></div> <div>80%</div> <div>15%</div> <div>..</div> </div> </div>

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	GOA	A	902	-	-	X	-
3	GOA	B	903	-	-	X	-
3	GOA	C	904	-	-	X	-
3	GOA	D	901	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6327 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 Acetate operon repressor.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	178	Total	C	N	O	S	Se	0	4	0
			1396	869	255	260	3	9			
1	B	179	Total	C	N	O	S	Se	0	7	0
			1424	892	258	260	3	11			
1	C	182	Total	C	N	O	S	Se	0	6	0
			1437	898	258	267	3	11			
1	D	178	Total	C	N	O	S	Se	0	5	0
			1407	881	256	256	3	11			

There are 48 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	GLY	-	cloning artifact	UNP P16528
A	2	HIS	-	cloning artifact	UNP P16528
A	3	MSE	-	cloning artifact	UNP P16528
A	19	MSE	MET	modified residue	UNP P16528
A	28	MSE	MET	modified residue	UNP P16528
A	50	MSE	MET	modified residue	UNP P16528
A	52	MSE	MET	modified residue	UNP P16528
A	62	MSE	MET	modified residue	UNP P16528
A	164	MSE	MET	modified residue	UNP P16528
A	179	MSE	MET	modified residue	UNP P16528
A	181	GLY	-	cloning artifact	UNP P16528
A	182	SER	-	cloning artifact	UNP P16528
B	1	GLY	-	cloning artifact	UNP P16528
B	2	HIS	-	cloning artifact	UNP P16528
B	3	MSE	-	cloning artifact	UNP P16528
B	19	MSE	MET	modified residue	UNP P16528
B	28	MSE	MET	modified residue	UNP P16528
B	50	MSE	MET	modified residue	UNP P16528
B	52	MSE	MET	modified residue	UNP P16528
B	62	MSE	MET	modified residue	UNP P16528
B	164	MSE	MET	modified residue	UNP P16528

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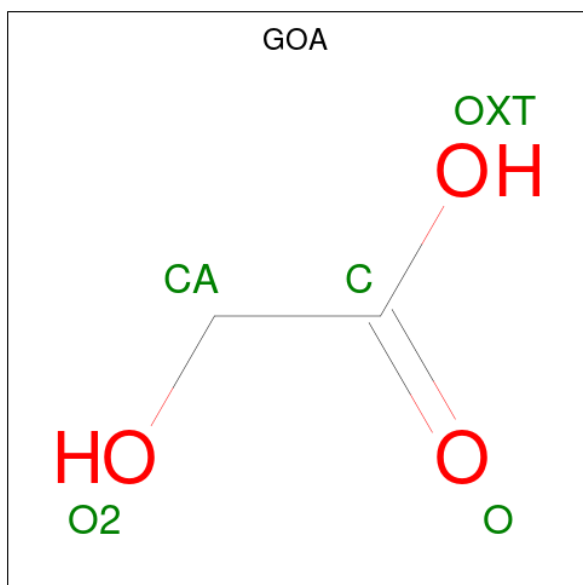
Chain	Residue	Modelled	Actual	Comment	Reference
B	179	MSE	MET	modified residue	UNP P16528
B	181	GLY	-	cloning artifact	UNP P16528
B	182	SER	-	cloning artifact	UNP P16528
C	1	GLY	-	cloning artifact	UNP P16528
C	2	HIS	-	cloning artifact	UNP P16528
C	3	MSE	-	cloning artifact	UNP P16528
C	19	MSE	MET	modified residue	UNP P16528
C	28	MSE	MET	modified residue	UNP P16528
C	50	MSE	MET	modified residue	UNP P16528
C	52	MSE	MET	modified residue	UNP P16528
C	62	MSE	MET	modified residue	UNP P16528
C	164	MSE	MET	modified residue	UNP P16528
C	179	MSE	MET	modified residue	UNP P16528
C	181	GLY	-	cloning artifact	UNP P16528
C	182	SER	-	cloning artifact	UNP P16528
D	1	GLY	-	cloning artifact	UNP P16528
D	2	HIS	-	cloning artifact	UNP P16528
D	3	MSE	-	cloning artifact	UNP P16528
D	19	MSE	MET	modified residue	UNP P16528
D	28	MSE	MET	modified residue	UNP P16528
D	50	MSE	MET	modified residue	UNP P16528
D	52	MSE	MET	modified residue	UNP P16528
D	62	MSE	MET	modified residue	UNP P16528
D	164	MSE	MET	modified residue	UNP P16528
D	179	MSE	MET	modified residue	UNP P16528
D	181	GLY	-	cloning artifact	UNP P16528
D	182	SER	-	cloning artifact	UNP P16528

- Molecule 2 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	B	1	Total	C	O	0	0
			4	2	2		
2	C	1	Total	C	O	0	0
			4	2	2		
2	D	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is GLYCOLIC ACID (three-letter code: GOA) (formula: $C_2H_4O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 5 2 3	0	0
3	B	1	Total C O 5 2 3	0	0
3	C	1	Total C O 5 2 3	0	0
3	D	1	Total C O 5 2 3	0	0

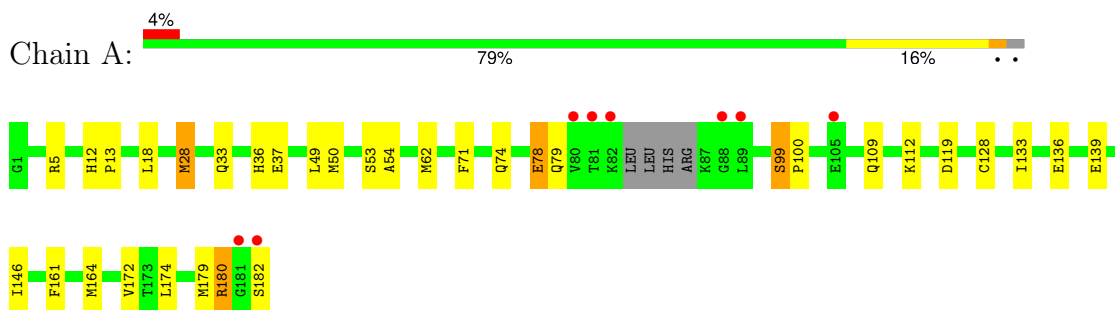
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	161	Total O 161 161	0	0
4	B	160	Total O 160 160	0	0
4	C	139	Total O 139 139	0	0
4	D	167	Total O 167 167	0	0

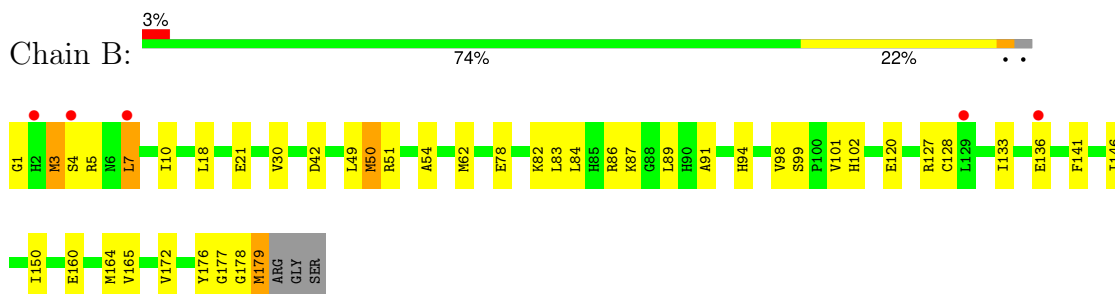
3 Residue-property plots [i](#)

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.

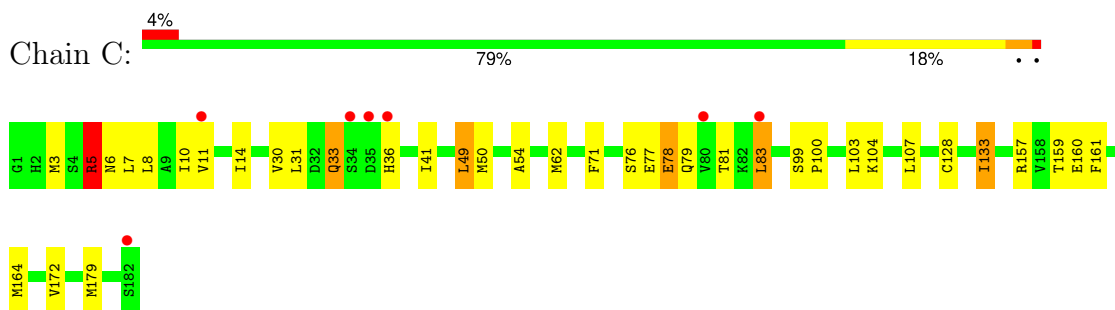
- Molecule 1: Acetate operon repressor



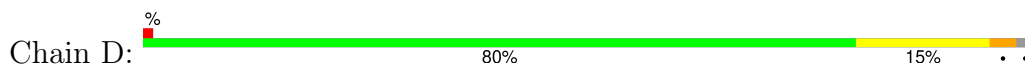
- Molecule 1: Acetate operon repressor

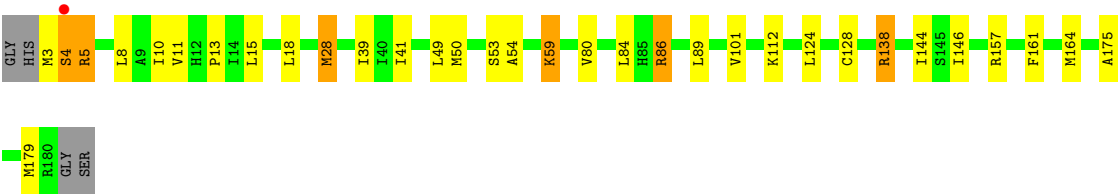


- Molecule 1: Acetate operon repressor



- Molecule 1: Acetate operon repressor





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.54Å 81.48Å 154.69Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 1.70 20.00 – 1.70	Depositor EDS
% Data completeness (in resolution range)	97.1 (20.00-1.70) 97.0 (20.00-1.70)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.32 (at 1.70Å)	Xtrriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.204 , 0.261 0.201 , 0.258	Depositor DCC
R_{free} test set	3677 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	25.3	Xtrriage
Anisotropy	0.015	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 44.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	6327	wwPDB-VP
Average B, all atoms (Å ²)	27.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 26.22 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 2.7405e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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: GOA, EDO

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	0.94	1/1410 (0.1%)	0.91	1/1882 (0.1%)
1	B	1.04	0/1449	0.93	4/1937 (0.2%)
1	C	1.05	0/1462	0.93	1/1953 (0.1%)
1	D	1.00	2/1432 (0.1%)	0.92	0/1915
All	All	1.01	3/5753 (0.1%)	0.92	6/7687 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	28[A]	MSE	CG-SE	-5.49	1.76	1.95
1	D	28[B]	MSE	CG-SE	-5.49	1.76	1.95
1	A	28	MSE	SE-CE	-5.26	1.64	1.95

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	5	ARG	NE-CZ-NH2	-6.11	117.25	120.30
1	B	3[A]	MSE	CB-CG-SE	-5.84	95.17	112.70
1	B	3[B]	MSE	CB-CG-SE	-5.84	95.17	112.70
1	A	119	ASP	CB-CG-OD2	-5.73	113.15	118.30
1	B	51	ARG	NE-CZ-NH2	-5.41	117.60	120.30
1	B	83	LEU	CB-CG-CD2	5.31	120.03	111.00

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	1396	0	1402	32	0
1	B	1424	0	1447	55	0
1	C	1437	0	1453	56	0
1	D	1407	0	1433	44	0
2	A	4	0	6	2	0
2	B	4	0	6	0	0
2	C	4	0	6	0	0
2	D	4	0	6	0	0
3	A	5	0	0	4	0
3	B	5	0	0	2	0
3	C	5	0	0	4	0
3	D	5	0	0	4	0
4	A	161	0	0	5	0
4	B	160	0	0	8	1
4	C	139	0	0	4	1
4	D	167	0	0	2	0
All	All	6327	0	5759	174	1

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 15.

All (174) 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:128:CYS:SG	3:B:903:GOA:CA	2.05	1.43
1:D:128:CYS:SG	3:D:901:GOA:CA	2.13	1.36
1:A:128:CYS:SG	3:A:902:GOA:CA	2.17	1.32
1:C:128:CYS:SG	3:C:904:GOA:CA	2.24	1.25
4:A:1036:HOH:O	1:B:3[A]:MSE:HB2	1.40	1.22
1:B:50[A]:MSE:HE1	1:C:50[A]:MSE:SE	1.98	1.12
1:A:161:PHE:HA	1:A:164:MSE:HE2	1.13	1.11
1:A:161:PHE:HA	1:A:164:MSE:CE	1.93	0.98
1:B:120:GLU:CG	1:B:150[B]:ILE:HD11	1.94	0.97
1:C:11:VAL:HG21	1:C:172:VAL:CG2	1.95	0.95
1:D:161:PHE:HA	1:D:164:MSE:HE3	1.49	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:128:CYS:HG	3:D:901:GOA:CA	1.76	0.94
1:B:120:GLU:HG2	1:B:150[B]:ILE:HD11	1.51	0.93
1:A:128:CYS:HG	3:A:902:GOA:CA	1.73	0.92
1:B:62[B]:MSE:SE	1:B:62[B]:MSE:H	2.02	0.92
1:B:7:LEU:HG	1:B:179:MSE:HE2	1.49	0.92
1:C:11:VAL:CG2	1:C:172:VAL:HG22	2.00	0.92
1:C:10:ILE:O	4:C:1038:HOH:O	1.88	0.91
1:A:179:MSE:HE3	1:A:182:SER:OG	1.72	0.90
1:C:11:VAL:HG21	1:C:172:VAL:HG22	1.56	0.87
1:C:3[B]:MSE:HE2	1:C:8:LEU:HG	1.56	0.86
1:C:3[B]:MSE:CA	1:C:3[B]:MSE:HE3	2.05	0.86
1:B:50[A]:MSE:CE	1:C:50[A]:MSE:SE	2.73	0.85
1:D:3:MSE:HE2	1:D:8:LEU:CD2	2.07	0.84
1:C:128:CYS:HG	3:C:904:GOA:CA	1.87	0.84
1:A:5[B]:ARG:HH22	1:B:5[B]:ARG:HH22	1.22	0.84
1:C:7:LEU:HA	1:C:179[B]:MSE:HE1	1.58	0.83
1:C:3[B]:MSE:HE3	1:C:3[B]:MSE:HA	1.63	0.81
1:B:3[B]:MSE:SE	1:B:7:LEU:HD22	2.30	0.81
1:C:11:VAL:HG23	1:C:14:ILE:HD12	1.65	0.78
1:D:15:LEU:HD11	1:D:28[A]:MSE:HE3	1.66	0.78
1:C:11:VAL:HG21	1:C:172:VAL:HG21	1.66	0.77
1:B:54:ALA:HB2	1:C:49:LEU:HD21	1.68	0.76
1:A:161:PHE:CA	1:A:164:MSE:HE2	2.07	0.76
1:D:86[A]:ARG:CG	1:D:86[A]:ARG:HH11	2.00	0.75
1:D:86[A]:ARG:HH11	1:D:86[A]:ARG:HG3	1.52	0.75
1:B:176:TYR:O	1:B:179:MSE:HG2	1.87	0.74
1:C:11:VAL:CG2	1:C:172:VAL:CG2	2.61	0.74
1:B:150[B]:ILE:HG22	4:B:955:HOH:O	1.88	0.73
1:D:28[A]:MSE:HE2	1:D:144:ILE:CG1	2.19	0.73
1:B:128:CYS:SG	3:B:903:GOA:C	2.76	0.73
1:C:7:LEU:CA	1:C:179[B]:MSE:HE1	2.18	0.73
1:D:84:LEU:HD22	1:D:89:LEU:HD21	1.72	0.72
1:B:21:GLU:OE2	4:B:986:HOH:O	2.09	0.71
1:A:50[B]:MSE:HE1	1:D:50[B]:MSE:HG2	1.73	0.70
1:C:128:CYS:SG	3:C:904:GOA:C	2.78	0.70
1:D:28[B]:MSE:SE	1:D:41:ILE:HD11	2.41	0.70
1:B:49:LEU:HD22	1:C:54:ALA:HB2	1.74	0.70
1:A:5[B]:ARG:NH2	1:B:5[B]:ARG:HH22	1.90	0.69
1:B:86[A]:ARG:NH1	4:B:1057:HOH:O	2.24	0.69
1:C:159:THR:HG22	4:C:966:HOH:O	1.93	0.69
1:D:128:CYS:SG	3:D:901:GOA:C	2.81	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:LEU:HG	1:B:179:MSE:CE	2.24	0.68
1:B:82:LYS:HG2	4:B:1056:HOH:O	1.93	0.67
1:B:120:GLU:CG	1:B:150[B]:ILE:CD1	2.71	0.67
1:B:18:LEU:HD11	1:B:146:ILE:HD11	1.78	0.65
1:D:39:ILE:HG12	1:D:59:LYS:HD3	1.76	0.65
1:A:128:CYS:SG	3:A:902:GOA:C	2.85	0.64
1:B:94:HIS:ND1	4:B:1023:HOH:O	2.26	0.64
1:C:11:VAL:HG23	1:C:172:VAL:HG22	1.77	0.64
1:A:12:HIS:HB3	1:A:13:PRO:HD3	1.80	0.64
1:C:3[A]:MSE:HG2	1:C:41:ILE:HD11	1.80	0.64
1:D:10:ILE:HD13	1:D:179[A]:MSE:SE	2.48	0.64
1:A:139:GLU:HG2	4:A:1055:HOH:O	1.99	0.62
1:C:7:LEU:HA	1:C:179[B]:MSE:CE	2.28	0.62
1:A:5[B]:ARG:HH12	1:B:5[B]:ARG:NH2	1.98	0.62
1:C:30:VAL:HG13	1:C:41:ILE:HD13	1.82	0.62
1:D:18:LEU:HD11	1:D:146:ILE:HD11	1.82	0.61
1:B:7:LEU:CG	1:B:179:MSE:HE2	2.27	0.61
1:D:3:MSE:CE	1:D:8:LEU:HG	2.31	0.61
1:D:3:MSE:HE3	1:D:4:SER:H	1.66	0.61
1:D:80:VAL:O	1:D:84:LEU:HG	2.02	0.60
1:D:3:MSE:HE2	1:D:8:LEU:HG	1.84	0.59
1:D:3:MSE:HE2	1:D:8:LEU:CG	2.33	0.59
1:C:157:ARG:O	1:C:160[A]:GLU:HB3	2.03	0.59
1:A:62:MSE:HE1	1:A:71:PHE:HB2	1.85	0.59
1:C:62:MSE:HB2	1:C:83:LEU:CD1	2.34	0.58
1:C:11:VAL:HG22	1:C:11:VAL:O	2.03	0.58
1:D:138:ARG:NH1	4:D:1057:HOH:O	2.34	0.58
1:B:84:LEU:HD22	1:B:89:LEU:HD21	1.85	0.58
1:A:74:GLN:NE2	2:A:801:EDO:H11	2.19	0.58
1:D:10:ILE:O	1:D:13:PRO:HD2	2.02	0.58
1:B:120:GLU:CD	1:B:150[B]:ILE:CD1	2.71	0.58
1:C:76:SER:H	1:C:79:GLN:HE21	1.53	0.57
1:D:86[A]:ARG:HG3	1:D:86[A]:ARG:NH1	2.14	0.57
1:D:28[A]:MSE:HE2	1:D:144:ILE:HD11	1.86	0.57
1:C:3[B]:MSE:CA	1:C:3[B]:MSE:CE	2.82	0.57
1:C:30:VAL:CG1	1:C:41:ILE:HD13	2.34	0.57
1:D:3:MSE:HG3	1:D:41:ILE:HD11	1.86	0.57
1:C:128:CYS:SG	3:C:904:GOA:O2	2.64	0.56
1:C:103:LEU:O	1:C:107:LEU:HG	2.06	0.56
1:B:62[B]:MSE:HE2	1:B:87:LYS:HE2	1.86	0.55
1:B:50[A]:MSE:SE	1:C:50[A]:MSE:SE	3.24	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:128:CYS:SG	3:D:901:GOA:O2	2.61	0.55
1:B:177:GLY:O	1:B:179:MSE:N	2.39	0.55
1:D:3:MSE:HE2	1:D:8:LEU:HD23	1.85	0.55
1:A:179:MSE:HE3	1:A:182:SER:HG	1.69	0.54
1:C:3[B]:MSE:CE	1:C:8:LEU:HG	2.33	0.54
1:A:78:GLU:HG3	1:A:79:GLN:N	2.22	0.54
1:B:136:GLU:OE1	1:B:176:TYR:HE2	1.91	0.53
1:D:28[A]:MSE:HE2	1:D:144:ILE:CD1	2.39	0.53
1:B:177:GLY:C	1:B:179:MSE:H	2.13	0.52
1:B:10:ILE:HD12	1:B:179:MSE:HE1	1.92	0.52
1:C:11:VAL:HG23	1:C:14:ILE:CD1	2.38	0.52
1:A:136:GLU:CD	1:A:180:ARG:HD3	2.30	0.52
1:B:3[A]:MSE:HG3	1:B:30:VAL:CG1	2.40	0.51
1:C:77:GLU:OE1	1:C:81:THR:OG1	2.29	0.51
1:C:33:GLN:NE2	1:C:33:GLN:H	2.08	0.51
1:B:4[A]:SER:O	1:B:5[A]:ARG:C	2.48	0.51
1:C:3[B]:MSE:HA	1:C:3[B]:MSE:CE	2.39	0.50
1:B:91:ALA:HB2	1:B:98:VAL:HG11	1.92	0.50
1:C:7:LEU:N	1:C:179[B]:MSE:HE1	2.25	0.50
1:C:161:PHE:HA	1:C:164:MSE:HE3	1.94	0.50
1:D:28[A]:MSE:HE2	1:D:144:ILE:HG12	1.92	0.50
1:B:1:GLY:N	4:B:994:HOH:O	2.27	0.50
1:A:128:CYS:SG	3:A:902:GOA:O2	2.63	0.49
1:D:175:ALA:O	1:D:179[A]:MSE:HG2	2.13	0.49
1:A:37:GLU:OE2	4:A:1002:HOH:O	2.19	0.49
1:A:36:HIS:O	4:A:1045:HOH:O	2.20	0.49
1:A:109:GLN:NE2	4:A:1047:HOH:O	2.45	0.49
1:B:3[B]:MSE:SE	1:B:7:LEU:HB3	2.63	0.49
1:D:4:SER:OG	1:D:5:ARG:N	2.46	0.48
1:D:28[A]:MSE:CE	1:D:144:ILE:HD11	2.44	0.48
1:D:28[A]:MSE:CE	1:D:144:ILE:HG12	2.43	0.48
1:A:5[B]:ARG:HD2	1:B:42:ASP:OD1	2.12	0.48
1:C:3[B]:MSE:HG3	1:C:41:ILE:HD11	1.96	0.48
1:C:76:SER:H	1:C:79:GLN:NE2	2.12	0.48
1:B:127:ARG:HG3	1:B:150[B]:ILE:HG12	1.95	0.48
1:A:74:GLN:HE21	2:A:801:EDO:H11	1.79	0.47
1:D:39:ILE:HG12	1:D:59:LYS:CD	2.44	0.47
1:B:99:SER:OG	1:B:101:VAL:HG12	2.15	0.47
1:A:18:LEU:HD11	1:A:146:ILE:HD11	1.96	0.47
1:C:62:MSE:HB2	1:C:83:LEU:HD11	1.97	0.47
1:B:160:GLU:HG2	1:B:164:MSE:HE3	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:5:ARG:HG3	1:C:6:ASN:N	2.30	0.47
1:C:133:ILE:HD13	1:C:133:ILE:N	2.29	0.47
1:D:161:PHE:HA	1:D:164:MSE:CE	2.33	0.47
1:B:4[A]:SER:O	1:B:7:LEU:N	2.48	0.46
1:A:49:LEU:HD22	1:D:54:ALA:HB2	1.96	0.46
1:D:3:MSE:HE3	1:D:4:SER:N	2.30	0.46
1:D:157[A]:ARG:HH11	1:D:157[A]:ARG:HG2	1.80	0.46
1:C:10:ILE:HD11	1:C:179[B]:MSE:HE3	1.96	0.45
1:C:62:MSE:HB2	1:C:83:LEU:HD13	1.97	0.45
1:B:86[B]:ARG:HG2	1:B:87:LYS:CD	2.46	0.45
1:B:133:ILE:HD13	1:B:172:VAL:HG11	1.98	0.45
1:B:82:LYS:HD2	1:B:86[A]:ARG:HH21	1.80	0.45
1:A:164:MSE:HE3	1:A:164:MSE:HB2	1.65	0.45
1:A:5[B]:ARG:NH1	1:B:5[B]:ARG:NH2	2.63	0.45
1:B:94:HIS:CD2	4:B:984:HOH:O	2.69	0.45
1:C:10:ILE:CD1	1:C:179[B]:MSE:HE3	2.47	0.44
1:B:3[A]:MSE:HG2	1:B:141:PHE:HE1	1.82	0.44
1:C:31:LEU:HD13	1:C:71:PHE:CG	2.53	0.44
1:B:101:VAL:HG13	1:B:102:HIS:N	2.32	0.44
1:C:100:PRO:O	1:C:104:LYS:HE2	2.18	0.44
1:C:49:LEU:HD23	1:C:49:LEU:HA	1.83	0.43
1:A:133:ILE:HD13	1:A:172:VAL:HG11	2.01	0.42
1:A:99:SER:HA	1:A:100:PRO:HD3	1.88	0.42
1:B:18:LEU:CD1	1:B:164:MSE:HB3	2.49	0.42
1:A:54:ALA:HB2	1:D:49:LEU:CD2	2.49	0.42
1:B:127:ARG:HD2	1:B:150[B]:ILE:HG23	2.01	0.41
1:C:5:ARG:NH2	4:C:941:HOH:O	2.53	0.41
1:C:160[A]:GLU:HG2	1:C:164:MSE:HE2	2.01	0.41
1:B:101:VAL:CG1	1:B:102:HIS:N	2.83	0.41
1:D:11:VAL:HG13	1:D:15:LEU:HD13	2.03	0.41
1:D:112:LYS:HE3	4:D:1031:HOH:O	2.19	0.41
1:D:28[B]:MSE:SE	1:D:41:ILE:CD1	3.17	0.40
1:D:3:MSE:HE2	1:D:8:LEU:HD21	1.97	0.40
1:C:78:GLU:OE2	4:C:919:HOH:O	2.22	0.40
1:A:174:LEU:HD23	1:A:174:LEU:HA	1.85	0.40
1:B:82:LYS:HD2	1:B:86[A]:ARG:NH2	2.36	0.40
1:B:78:GLU:HB2	4:B:948:HOH:O	2.22	0.40
1:C:31:LEU:HD13	1:C:71:PHE:CD1	2.56	0.40
1:D:86[A]:ARG:CG	1:D:86[A]:ARG:NH1	2.67	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:973:HOH:O	4:C:1041:HOH:O[1_455]	2.11	0.09

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	178/182 (98%)	172 (97%)	5 (3%)	1 (1%)	22	10
1	B	184/182 (101%)	178 (97%)	5 (3%)	1 (0%)	25	12
1	C	186/182 (102%)	179 (96%)	7 (4%)	0	100	100
1	D	181/182 (100%)	173 (96%)	7 (4%)	1 (1%)	22	10
All	All	729/728 (100%)	702 (96%)	24 (3%)	3 (0%)	30	17

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	180	ARG
1	B	178	GLY
1	D	5	ARG

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	148/140 (106%)	142 (96%)	6 (4%)	26	10
1	B	153/140 (109%)	148 (97%)	5 (3%)	33	16

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	C	154/140 (110%)	146 (95%)	8 (5%)	19 7
1	D	151/140 (108%)	143 (95%)	8 (5%)	19 6
All	All	606/560 (108%)	579 (96%)	27 (4%)	25 9

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	28	MSE
1	A	33	GLN
1	A	53	SER
1	A	78	GLU
1	A	99	SER
1	A	112	LYS
1	B	7	LEU
1	B	50[A]	MSE
1	B	50[B]	MSE
1	B	165	VAL
1	B	179	MSE
1	C	5	ARG
1	C	33	GLN
1	C	36	HIS
1	C	49	LEU
1	C	78	GLU
1	C	83	LEU
1	C	99	SER
1	C	133	ILE
1	D	4	SER
1	D	53	SER
1	D	59	LYS
1	D	86[A]	ARG
1	D	86[B]	ARG
1	D	101	VAL
1	D	124	LEU
1	D	138	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	33	GLN
1	A	48	HIS

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Mol	Chain	Res	Type
1	A	74	GLN
1	A	109	GLN
1	B	6	ASN
1	B	33	GLN
1	C	33	GLN
1	C	36	HIS
1	C	74	GLN
1	C	79	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

8 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	EDO	C	803	-	3,3,3	0.46	0	2,2,2	0.30	0
3	GOA	B	903	-	4,4,4	1.09	0	4,4,4	1.38	1 (25%)
3	GOA	D	901	-	4,4,4	1.48	1 (25%)	4,4,4	1.87	1 (25%)
3	GOA	A	902	-	4,4,4	1.09	0	4,4,4	2.18	1 (25%)
2	EDO	B	802	-	3,3,3	0.59	0	2,2,2	0.54	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	EDO	A	801	-	3,3,3	0.46	0	2,2,2	1.00	0
2	EDO	D	804	-	3,3,3	0.43	0	2,2,2	0.19	0
3	GOA	C	904	-	4,4,4	1.26	1 (25%)	4,4,4	1.82	1 (25%)

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	EDO	C	803	-	-	0/1/1/1	-
3	GOA	B	903	-	-	0/2/2/2	-
3	GOA	D	901	-	-	0/2/2/2	-
3	GOA	A	902	-	-	0/2/2/2	-
2	EDO	B	802	-	-	1/1/1/1	-
2	EDO	A	801	-	-	1/1/1/1	-
2	EDO	D	804	-	-	0/1/1/1	-
3	GOA	C	904	-	-	0/2/2/2	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	D	901	GOA	OXT-C	-2.41	1.22	1.30
3	C	904	GOA	OXT-C	-2.32	1.23	1.30

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	902	GOA	OXT-C-CA	4.01	119.19	111.90
3	D	901	GOA	OXT-C-CA	3.40	118.07	111.90
3	C	904	GOA	OXT-C-CA	3.38	118.04	111.90
3	B	903	GOA	OXT-C-CA	2.35	116.17	111.90

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	801	EDO	O1-C1-C2-O2
2	B	802	EDO	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	903	GOA	2	0
3	D	901	GOA	4	0
3	A	902	GOA	4	0
2	A	801	EDO	2	0
3	C	904	GOA	4	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	170/182 (93%)	0.25	8 (4%) 37 40	9, 27, 45, 63	3 (1%)
1	B	171/182 (93%)	0.05	5 (2%) 54 57	9, 22, 41, 56	6 (3%)
1	C	174/182 (95%)	0.05	7 (4%) 43 46	10, 22, 37, 51	3 (1%)
1	D	170/182 (93%)	-0.03	1 (0%) 85 88	13, 23, 39, 52	2 (1%)
All	All	685/728 (94%)	0.08	21 (3%) 51 54	9, 24, 42, 63	14 (2%)

All (21) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	34	SER	5.2
1	C	80	VAL	3.7
1	A	89	LEU	3.0
1	C	11	VAL	2.7
1	A	88	GLY	2.7
1	C	83	LEU	2.7
1	C	35	ASP	2.7
1	D	4	SER	2.5
1	B	2	HIS	2.4
1	B	129	LEU	2.4
1	B	136	GLU	2.3
1	C	36	HIS	2.3
1	A	182	SER	2.3
1	A	81	THR	2.3
1	A	105	GLU	2.2
1	B	4[A]	SER	2.2
1	A	181	GLY	2.2
1	A	80	VAL	2.1
1	B	7	LEU	2.1
1	A	82	LYS	2.0
1	C	182	SER	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, 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
2	EDO	A	801	4/4	0.88	0.10	32,33,35,36	0
2	EDO	C	803	4/4	0.89	0.12	23,27,29,30	0
2	EDO	B	802	4/4	0.90	0.10	26,31,34,38	0
3	GOA	A	902	5/5	0.94	0.07	22,24,27,28	0
3	GOA	D	901	5/5	0.95	0.09	18,21,22,23	0
3	GOA	B	903	5/5	0.96	0.05	18,18,20,22	0
2	EDO	D	804	4/4	0.96	0.05	22,23,24,24	0
3	GOA	C	904	5/5	0.97	0.04	20,20,23,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.