



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 03:02 PM UTC

PDB ID : 9D5W / pdb_00009d5w
Title : Crystal Structure of the Substrate Binding Domain Protein of the ABC Transporter PBP2_YxeM from *Vibrio cholerae*
Authors : Kim, Y.; Maltseva, N.; Grimshaw, S.; Joachimiak, A.; Center for Structural Biology of Infectious Diseases (CSBID)
Deposited on : 2024-08-14
Resolution : 2.30 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

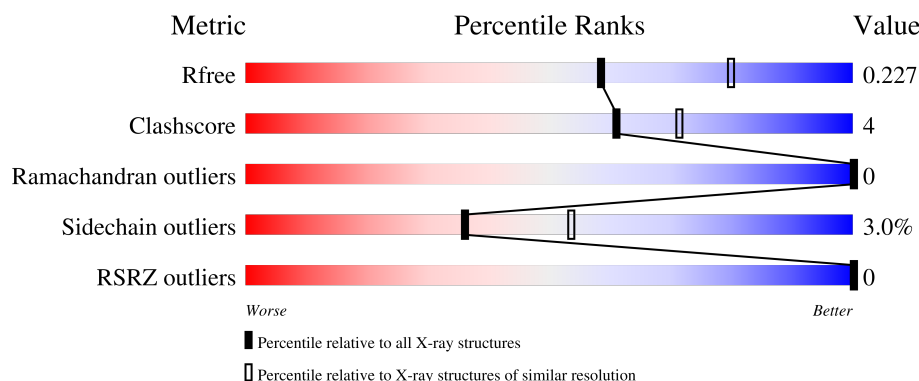
1 Overall quality at a glance ⓘ

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	233	 79% 18% . .
1	B	233	 81% 16% . .
1	C	233	 82% 15% .
1	D	233	 86% 11% .
1	E	233	 85% 13% .

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Mol	Chain	Length	Quality of chain
1	F	233	 89% 8% •
1	G	233	 87% 11% •
1	H	233	 84% 13% •
1	I	233	 85% 12% •

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
6	SO4	G	303	-	-	X	-

2 Entry composition

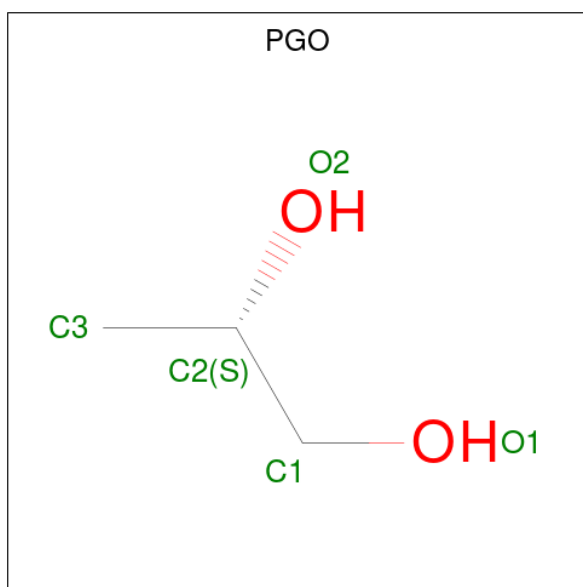
There are 7 unique types of molecules in this entry. The entry contains 16666 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Amino acid ABC transporter, periplasmic amino acid-binding portion.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	227	Total	C	N	O	S	0	0	0
			1770	1121	301	345	3			
1	B	227	Total	C	N	O	S	0	0	0
			1770	1121	301	345	3			
1	C	227	Total	C	N	O	S	0	0	0
			1770	1121	301	345	3			
1	D	227	Total	C	N	O	S	0	0	0
			1770	1121	301	345	3			
1	E	227	Total	C	N	O	S	0	0	0
			1770	1121	301	345	3			
1	F	227	Total	C	N	O	S	0	0	0
			1770	1121	301	345	3			
1	G	227	Total	C	N	O	S	0	0	0
			1770	1121	301	345	3			
1	H	227	Total	C	N	O	S	0	0	0
			1770	1121	301	345	3			
1	I	227	Total	C	N	O	S	0	0	0
			1770	1121	301	345	3			

- Molecule 2 is S-1,2-PROPANEDIOL (CCD ID: PGO) (formula: C₃H₈O₂).



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

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Cl	0	0
			2	2		
3	B	2	Total	Cl	0	0
			2	2		
3	C	2	Total	Cl	0	0
			2	2		

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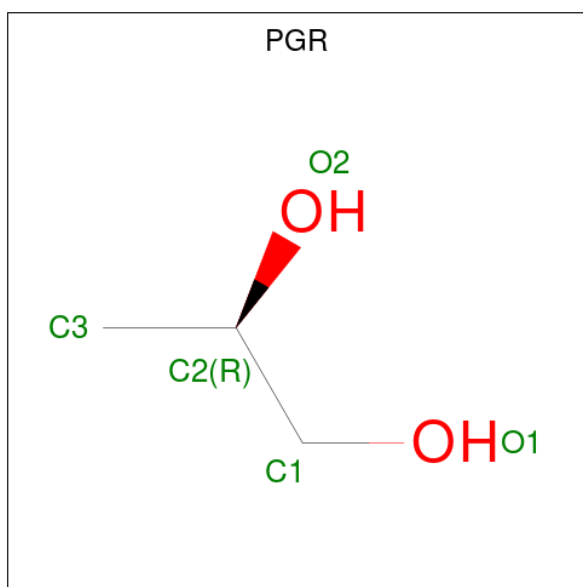
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	1	Total 1	Cl 1	0	0
3	F	4	Total 4	Cl 4	0	0
3	G	2	Total 2	Cl 2	0	0
3	H	2	Total 2	Cl 2	0	0
3	I	2	Total 2	Cl 2	0	0

- Molecule 4 is POTASSIUM ION (CCD ID: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	1	Total 1	K 1	0	0
4	B	1	Total 1	K 1	0	0
4	C	1	Total 1	K 1	0	0
4	D	1	Total 1	K 1	0	0
4	E	1	Total 1	K 1	0	0
4	G	1	Total 1	K 1	0	0

- Molecule 5 is R-1,2-PROPANEDIOL (CCD ID: PGR) (formula: C₃H₈O₂).



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

- Molecule 6 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	D	1	Total	O	S	0	0
			5	4	1		
6	G	1	Total	O	S	0	0
			5	4	1		
6	G	1	Total	O	S	0	0
			5	4	1		

- Molecule 7 is water.

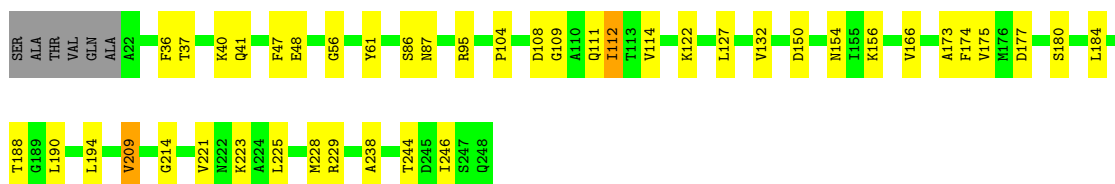
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	76	Total 76	O 76	0	0
7	B	70	Total 70	O 70	0	0
7	C	39	Total 39	O 39	0	0
7	D	83	Total 83	O 83	0	0
7	E	88	Total 88	O 88	0	0
7	F	74	Total 74	O 74	0	0
7	G	47	Total 47	O 47	0	0
7	H	83	Total 83	O 83	0	0
7	I	88	Total 88	O 88	0	0

3 Residue-property plots

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.

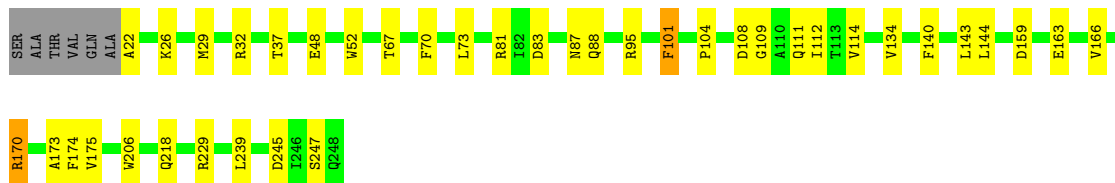
- Molecule 1: Amino acid ABC transporter, periplasmic amino acid-binding portion

Chain A: 



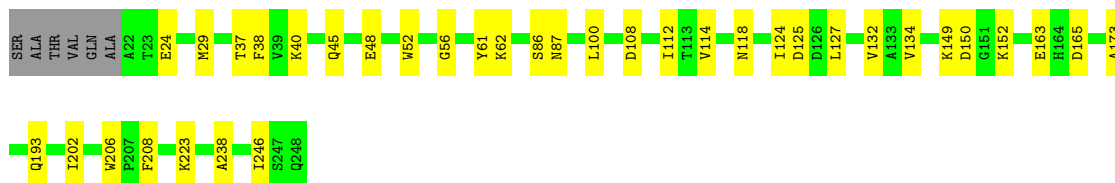
- Molecule 1: Amino acid ABC transporter, periplasmic amino acid-binding portion

Chain B: 



- Molecule 1: Amino acid ABC transporter, periplasmic amino acid-binding portion

Chain C: 



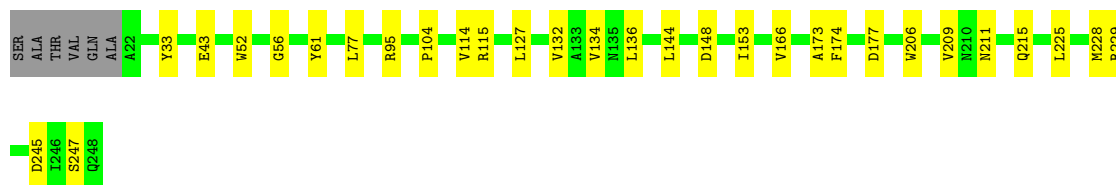
- Molecule 1: Amino acid ABC transporter, periplasmic amino acid-binding portion

Chain D: 



- Molecule 1: Amino acid ABC transporter, periplasmic amino acid-binding portion

Chain E:  85% 13% .



- Molecule 1: Amino acid ABC transporter, periplasmic amino acid-binding portion

Chain F:  89% 8% .



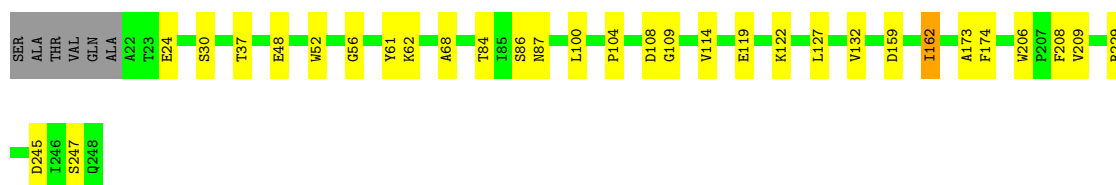
- Molecule 1: Amino acid ABC transporter, periplasmic amino acid-binding portion

Chain G:  87% 11% .



- Molecule 1: Amino acid ABC transporter, periplasmic amino acid-binding portion

Chain H:  84% 13% .



- Molecule 1: Amino acid ABC transporter, periplasmic amino acid-binding portion

Chain I:  85% 12% .



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	118.04Å 105.82Å 118.12Å 90.00° 119.92° 90.00°	Depositor
Resolution (Å)	59.11 – 2.30 59.11 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.5 (59.11-2.30) 95.5 (59.11-2.30)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.05 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.8.0430	Depositor
R, R_{free}	0.194 , 0.229 0.193 , 0.227	Depositor DCC
R_{free} test set	5579 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	28.4	Xtriage
Anisotropy	0.859	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 17.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.39$, $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	0.429 for -h-l,k,h 0.429 for l,k,-h-l 0.188 for h,-k,-h-l 0.190 for -h-l,-k,l 0.188 for l,-k,h	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	16666	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.54% 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: CL, SO4, PGR, K, PGO

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	0.09	0/1799	0.24	0/2429
1	B	0.09	0/1799	0.23	0/2429
1	C	0.09	0/1799	0.23	0/2429
1	D	0.09	0/1799	0.24	0/2429
1	E	0.09	0/1799	0.24	0/2429
1	F	0.09	0/1799	0.24	0/2429
1	G	0.09	0/1799	0.23	0/2429
1	H	0.09	0/1799	0.25	0/2429
1	I	0.09	0/1799	0.23	0/2429
All	All	0.09	0/16191	0.24	0/21861

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

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	1770	0	1757	20	0
1	B	1770	0	1756	24	0
1	C	1770	0	1756	16	0
1	D	1770	0	1757	12	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	1770	0	1757	16	0
1	F	1770	0	1757	9	0
1	G	1770	0	1757	13	0
1	H	1770	0	1757	14	0
1	I	1770	0	1757	16	0
2	A	10	0	16	0	0
2	B	10	0	16	0	0
2	D	15	0	24	0	0
2	G	5	0	8	0	0
2	I	5	0	8	0	0
3	A	2	0	0	1	0
3	B	2	0	0	1	0
3	C	2	0	0	0	0
3	E	1	0	0	1	0
3	F	4	0	0	0	0
3	G	2	0	0	1	0
3	H	2	0	0	0	0
3	I	2	0	0	0	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
4	E	1	0	0	0	0
4	G	1	0	0	0	0
5	C	5	0	8	0	0
6	D	5	0	0	0	0
6	G	10	0	0	3	0
7	A	76	0	0	0	0
7	B	70	0	0	2	0
7	C	39	0	0	1	0
7	D	83	0	0	1	0
7	E	88	0	0	0	0
7	F	74	0	0	1	0
7	G	47	0	0	2	0
7	H	83	0	0	0	0
7	I	88	0	0	2	0
All	All	16666	0	15891	135	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 4.

All (135) 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:95:ARG:NH1	3:A:304:CL:CL	2.63	0.65
1:D:29:MET:HB2	1:D:37:THR:HG21	1.78	0.63
1:A:112:ILE:HD13	1:A:175:VAL:HG22	1.80	0.63
1:E:56:GLY:HA2	1:E:61:TYR:HB2	1.85	0.59
1:G:162:ILE:HD11	1:G:176:MET:HE3	1.83	0.59
1:D:31:GLY:HA2	1:D:37:THR:HG22	1.85	0.59
1:E:104:PRO:O	1:E:229:ARG:NH2	2.37	0.57
1:G:170:ARG:NH1	6:G:303:SO4:O1	2.37	0.57
1:F:115:ARG:NH1	7:F:403:HOH:O	2.38	0.56
1:I:158:TYR:HB2	1:I:162:ILE:HG22	1.87	0.56
1:B:48:GLU:OE2	1:B:88:GLN:NE2	2.35	0.56
1:H:119:GLU:O	1:H:122:LYS:NZ	2.38	0.55
1:D:114:VAL:HG12	1:D:173:ALA:HB2	1.88	0.55
1:B:81:ARG:HD3	1:I:136:LEU:HD21	1.90	0.54
1:C:124:ILE:HD12	1:C:127:LEU:HD12	1.91	0.53
1:B:101:PHE:O	1:B:218:GLN:NE2	2.41	0.53
1:C:48:GLU:OE1	1:C:86:SER:OG	2.23	0.53
1:E:77:LEU:HG	1:E:209:VAL:HG13	1.91	0.52
1:B:95:ARG:NH1	3:B:303:CL:CL	2.74	0.52
1:A:114:VAL:HG12	1:A:173:ALA:HB2	1.92	0.52
1:H:104:PRO:O	1:H:229:ARG:NH2	2.40	0.51
1:H:24:GLU:HA	1:H:62:LYS:HB3	1.92	0.51
1:H:127:LEU:HD13	1:H:132:VAL:HG21	1.92	0.51
1:A:127:LEU:HD13	1:A:132:VAL:HG21	1.93	0.51
1:H:48:GLU:OE1	1:H:86:SER:OG	2.28	0.51
1:I:127:LEU:HD13	1:I:132:VAL:HG21	1.93	0.51
1:B:140:PHE:HA	1:B:143:LEU:HD12	1.94	0.50
1:B:22:ALA:N	7:B:402:HOH:O	2.44	0.50
1:C:24:GLU:HA	1:C:62:LYS:HB3	1.93	0.50
1:G:170:ARG:NH1	6:G:303:SO4:S	2.84	0.50
1:E:225:LEU:HA	1:E:228:MET:HE2	1.93	0.50
1:A:108:ASP:CG	1:A:109:GLY:H	2.20	0.49
1:B:114:VAL:HG12	1:B:173:ALA:HB2	1.94	0.49
1:H:114:VAL:HG12	1:H:173:ALA:HB2	1.94	0.49
1:B:159:ASP:OD2	1:I:32:ARG:NH2	2.45	0.49
1:F:136:LEU:N	1:F:158:TYR:O	2.39	0.49
1:A:104:PRO:O	1:A:229:ARG:NH2	2.46	0.48
1:G:114:VAL:HG12	1:G:173:ALA:HB2	1.94	0.48
1:G:111:GLN:HG2	1:G:194:LEU:HD22	1.95	0.48
1:B:166:VAL:HG21	1:B:174:PHE:HD1	1.79	0.47
1:A:154:ASN:OD1	1:A:156:LYS:NZ	2.47	0.47
1:C:114:VAL:HG12	1:C:173:ALA:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:30:SER:HA	1:H:68:ALA:O	2.14	0.47
1:I:22:ALA:N	7:I:407:HOH:O	2.48	0.47
1:I:92:THR:O	1:I:96:LYS:N	2.46	0.47
1:A:177:ASP:OD1	1:A:180:SER:OG	2.31	0.47
1:B:111:GLN:NE2	1:B:112:ILE:O	2.48	0.47
1:D:32:ARG:NH2	7:D:409:HOH:O	2.45	0.47
1:D:112:ILE:HD11	1:D:124:ILE:HD12	1.98	0.46
1:C:118:ASN:O	1:C:193:GLN:NE2	2.48	0.46
1:E:52:TRP:HE1	1:E:206:TRP:CD1	2.33	0.46
1:B:104:PRO:O	1:B:229:ARG:NH2	2.45	0.46
1:G:29:MET:HB2	1:G:37:THR:HG21	1.97	0.46
1:B:144:LEU:HD21	1:B:175:VAL:HG21	1.97	0.46
1:D:48:GLU:OE1	1:D:86:SER:OG	2.32	0.46
1:D:126:ASP:N	1:D:126:ASP:OD1	2.49	0.46
1:I:182:LEU:HD21	1:I:194:LEU:HD11	1.98	0.46
1:C:52:TRP:HE1	1:C:206:TRP:CD1	2.34	0.46
1:B:29:MET:HB2	1:B:37:THR:HG21	1.97	0.45
1:C:223:LYS:HE2	1:C:223:LYS:HB3	1.83	0.45
1:F:108:ASP:CG	1:F:109:GLY:H	2.24	0.45
1:I:184:LEU:O	1:I:188:THR:OG1	2.31	0.45
1:F:167:ALA:HB1	1:F:191:PRO:HG2	1.98	0.45
1:B:32:ARG:NH2	7:B:403:HOH:O	2.49	0.45
1:H:162:ILE:HD11	1:H:174:PHE:HB2	1.97	0.45
1:D:27:VAL:HG13	1:D:84:THR:HG23	1.97	0.45
1:B:52:TRP:HE1	1:B:206:TRP:NE1	2.15	0.45
1:A:223:LYS:HA	1:B:239:LEU:HD13	1.98	0.44
3:G:305:CL:CL	7:G:436:HOH:O	2.58	0.44
1:C:238:ALA:HB1	1:C:246:ILE:HG12	1.99	0.44
1:A:56:GLY:HA2	1:A:61:TYR:HB2	2.00	0.44
1:A:111:GLN:HG2	1:A:194:LEU:HD22	2.00	0.44
1:E:114:VAL:HG12	1:E:173:ALA:HB2	2.00	0.44
1:B:112:ILE:HG12	1:B:175:VAL:HG22	1.99	0.44
1:C:29:MET:HB2	1:C:37:THR:HG21	1.98	0.44
1:H:52:TRP:CZ2	1:H:84:THR:HG21	2.53	0.43
1:H:108:ASP:CG	1:H:109:GLY:H	2.26	0.43
1:D:111:GLN:N	1:D:176:MET:O	2.45	0.43
1:E:136:LEU:HD22	1:G:168:LEU:HD23	2.00	0.43
1:E:148:ASP:HB2	1:E:153:ILE:HD12	2.00	0.43
1:F:144:LEU:HD21	1:F:175:VAL:HG21	2.00	0.43
1:D:56:GLY:HA2	1:D:61:TYR:HB2	2.01	0.43
1:E:166:VAL:HG21	1:E:174:PHE:HD2	1.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:28:GLY:HA2	1:I:66:VAL:O	2.19	0.43
1:A:36:PHE:HA	1:A:47:PHE:HB3	2.00	0.43
1:C:127:LEU:HD13	1:C:132:VAL:HG21	2.00	0.43
1:A:184:LEU:HB3	1:A:190:LEU:HD12	2.00	0.43
1:A:225:LEU:HA	1:A:228:MET:HE2	2.00	0.43
1:H:56:GLY:HA2	1:H:61:TYR:HB2	2.00	0.43
1:C:100:LEU:O	1:C:208:PHE:N	2.50	0.43
1:G:30:SER:HA	1:G:68:ALA:O	2.19	0.42
1:H:100:LEU:O	1:H:208:PHE:N	2.52	0.42
1:G:31:GLY:C	1:G:32:ARG:HE	2.27	0.42
1:D:112:ILE:HG12	1:D:175:VAL:HG22	2.00	0.42
1:E:95:ARG:NH2	3:E:301:CL:CL	2.82	0.42
1:I:84:THR:OG1	1:I:85:ILE:N	2.52	0.42
1:A:48:GLU:OE1	1:A:86:SER:OG	2.26	0.42
1:F:115:ARG:HG3	1:F:167:ALA:HA	2.02	0.42
1:C:56:GLY:HA2	1:C:61:TYR:HB2	2.02	0.42
1:I:163:GLU:H	1:I:163:GLU:CD	2.26	0.42
1:B:70:PHE:HA	1:B:73:LEU:HG	2.01	0.42
1:A:188:THR:HA	1:E:115:ARG:HH21	1.85	0.42
1:C:152:LYS:NZ	7:C:401:HOH:O	2.53	0.42
1:I:103:ASP:HB2	1:I:225:LEU:HD23	2.02	0.42
1:E:245:ASP:OD1	1:E:247:SER:OG	2.38	0.42
1:H:245:ASP:OD1	1:H:247:SER:OG	2.31	0.42
1:A:209:VAL:O	1:A:214:GLY:HA3	2.19	0.41
1:G:24:GLU:HG2	1:G:62:LYS:HD3	2.01	0.41
1:I:149:LYS:N	7:I:412:HOH:O	2.52	0.41
1:E:33:TYR:OH	1:E:177:ASP:OD1	2.28	0.41
1:A:238:ALA:HB1	1:A:246:ILE:HG12	2.03	0.41
1:C:108:ASP:HB3	1:C:202:ILE:HD11	2.01	0.41
1:G:34:PHE:CZ	1:G:187:LYS:HD2	2.55	0.41
1:H:52:TRP:HE1	1:H:206:TRP:CD1	2.38	0.41
1:B:163:GLU:H	1:B:163:GLU:CD	2.28	0.41
1:B:170:ARG:NH2	1:C:38:PHE:HB2	2.36	0.41
1:F:103:ASP:HB2	1:F:225:LEU:HD23	2.03	0.41
1:I:104:PRO:O	1:I:229:ARG:NH2	2.54	0.41
1:G:223:LYS:NZ	7:G:403:HOH:O	2.53	0.41
1:A:40:LYS:HG2	1:A:41:GLN:HG3	2.01	0.41
1:B:29:MET:O	1:B:67:THR:HA	2.21	0.41
1:B:245:ASP:OD1	1:B:247:SER:OG	2.38	0.41
1:D:108:ASP:CG	1:D:109:GLY:H	2.28	0.41
1:F:88:GLN:OE1	1:F:204:ASN:ND2	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:98:LYS:HE3	1:I:98:LYS:HB2	1.86	0.41
1:G:170:ARG:NH1	6:G:303:SO4:O3	2.54	0.41
1:I:166:VAL:HG22	1:I:171:ALA:HB3	2.02	0.40
1:B:26:LYS:O	1:B:83:ASP:N	2.49	0.40
1:F:58:ARG:NH2	1:F:231:ASP:OD2	2.52	0.40
1:E:52:TRP:HE1	1:E:206:TRP:NE1	2.19	0.40
1:E:211:ASN:O	1:E:215:GLN:N	2.46	0.40
1:A:166:VAL:HG21	1:A:174:PHE:HD2	1.87	0.40
1:B:108:ASP:CG	1:B:109:GLY:H	2.29	0.40
1:C:40:LYS:HB3	1:C:45:GLN:NE2	2.37	0.40
1:E:127:LEU:HD13	1:E:132:VAL:HG21	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	225/233 (97%)	216 (96%)	9 (4%)	0	100	100
1	B	225/233 (97%)	211 (94%)	14 (6%)	0	100	100
1	C	225/233 (97%)	219 (97%)	6 (3%)	0	100	100
1	D	225/233 (97%)	213 (95%)	12 (5%)	0	100	100
1	E	225/233 (97%)	214 (95%)	11 (5%)	0	100	100
1	F	225/233 (97%)	212 (94%)	13 (6%)	0	100	100
1	G	225/233 (97%)	216 (96%)	9 (4%)	0	100	100
1	H	225/233 (97%)	212 (94%)	13 (6%)	0	100	100
1	I	225/233 (97%)	217 (96%)	8 (4%)	0	100	100
All	All	2025/2097 (97%)	1930 (95%)	95 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	187/191 (98%)	179 (96%)	8 (4%)	26	39
1	B	187/191 (98%)	183 (98%)	4 (2%)	47	66
1	C	187/191 (98%)	179 (96%)	8 (4%)	26	39
1	D	187/191 (98%)	179 (96%)	8 (4%)	26	39
1	E	187/191 (98%)	184 (98%)	3 (2%)	55	73
1	F	187/191 (98%)	182 (97%)	5 (3%)	39	58
1	G	187/191 (98%)	181 (97%)	6 (3%)	34	51
1	H	187/191 (98%)	182 (97%)	5 (3%)	39	58
1	I	187/191 (98%)	184 (98%)	3 (2%)	55	73
All	All	1683/1719 (98%)	1633 (97%)	50 (3%)	36	53

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	37	THR
1	A	87	ASN
1	A	112	ILE
1	A	122	LYS
1	A	150	ASP
1	A	209	VAL
1	A	221	VAL
1	A	244	THR
1	B	87	ASN
1	B	101	PHE
1	B	134	VAL
1	B	170	ARG
1	C	87	ASN
1	C	112	ILE
1	C	125	ASP
1	C	134	VAL
1	C	149	LYS
1	C	150	ASP

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Mol	Chain	Res	Type
1	C	163	GLU
1	C	165	ASP
1	D	87	ASN
1	D	101	PHE
1	D	126	ASP
1	D	134	VAL
1	D	144	LEU
1	D	162	ILE
1	D	170	ARG
1	D	179	LEU
1	E	43	GLU
1	E	134	VAL
1	E	144	LEU
1	F	40	LYS
1	F	125	ASP
1	F	144	LEU
1	F	162	ILE
1	F	163	GLU
1	G	134	VAL
1	G	156	LYS
1	G	160	THR
1	G	163	GLU
1	G	184	LEU
1	G	217	LEU
1	H	37	THR
1	H	87	ASN
1	H	159	ASP
1	H	162	ILE
1	H	209	VAL
1	I	126	ASP
1	I	135	ASN
1	I	144	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (14) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	142	GLN
1	A	218	GLN
1	B	135	ASN
1	C	118	ASN
1	D	142	GLN
1	E	204	ASN

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Mol	Chain	Res	Type
1	F	69	ASN
1	F	87	ASN
1	G	45	GLN
1	G	87	ASN
1	H	69	ASN
1	H	142	GLN
1	I	41	GLN
1	I	154	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 36 ligands modelled in this entry, 23 are monoatomic - leaving 13 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	$\# Z > 2$
2	PGO	B	302	-	4,4,4	0.56	0	4,4,4	0.88	0
2	PGO	D	302	-	4,4,4	0.57	0	4,4,4	0.88	0
2	PGO	G	301	-	4,4,4	0.61	0	4,4,4	0.85	0
6	SO4	D	304	-	4,4,4	0.24	0	6,6,6	0.09	0
2	PGO	I	301	-	4,4,4	0.58	0	4,4,4	0.90	0
6	SO4	G	303	-	4,4,4	0.23	0	6,6,6	0.07	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
6	SO4	G	302	-	4,4,4	0.24	0	6,6,6	0.08	0
2	PGO	D	301	-	4,4,4	0.57	0	4,4,4	0.91	0
2	PGO	B	301	-	4,4,4	0.58	0	4,4,4	0.89	0
2	PGO	A	302	-	4,4,4	0.57	0	4,4,4	0.90	0
2	PGO	A	301	-	4,4,4	0.58	0	4,4,4	0.90	0
2	PGO	D	303	-	4,4,4	0.58	0	4,4,4	0.92	0
5	PGR	C	301	-	4,4,4	0.63	0	4,4,4	0.62	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	PGO	B	302	-	-	0/2/2/2	-
2	PGO	D	302	-	-	0/2/2/2	-
2	PGO	G	301	-	-	0/2/2/2	-
2	PGO	I	301	-	-	0/2/2/2	-
2	PGO	D	301	-	-	2/2/2/2	-
2	PGO	B	301	-	-	0/2/2/2	-
2	PGO	A	302	-	-	2/2/2/2	-
2	PGO	A	301	-	-	2/2/2/2	-
2	PGO	D	303	-	-	2/2/2/2	-
5	PGR	C	301	-	-	1/2/2/2	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (9) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	301	PGO	O1-C1-C2-C3
2	A	301	PGO	O1-C1-C2-O2
2	A	302	PGO	O1-C1-C2-O2
2	D	301	PGO	O1-C1-C2-C3
2	D	301	PGO	O1-C1-C2-O2
2	D	303	PGO	O1-C1-C2-C3
2	A	302	PGO	O1-C1-C2-C3
2	D	303	PGO	O1-C1-C2-O2
5	C	301	PGR	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	G	303	SO4	3	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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	227/233 (97%)	-1.33	0 100 100	20, 29, 48, 65	0
1	B	227/233 (97%)	-1.39	0 100 100	20, 28, 41, 48	0
1	C	227/233 (97%)	-1.35	0 100 100	21, 28, 45, 61	0
1	D	227/233 (97%)	-1.38	0 100 100	17, 27, 41, 57	0
1	E	227/233 (97%)	-1.36	0 100 100	21, 28, 42, 61	0
1	F	227/233 (97%)	-1.38	0 100 100	20, 27, 40, 47	0
1	G	227/233 (97%)	-1.35	0 100 100	21, 28, 39, 54	0
1	H	227/233 (97%)	-1.36	0 100 100	20, 28, 45, 57	0
1	I	227/233 (97%)	-1.38	0 100 100	19, 27, 39, 68	0
All	All	2043/2097 (97%)	-1.37	0 100 100	17, 28, 43, 68	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	PGO	A	302	5/5	0.97	0.06	32,34,35,36	0
2	PGO	D	303	5/5	0.97	0.06	35,35,35,36	0
2	PGO	A	301	5/5	0.98	0.05	24,25,26,26	0
3	CL	F	304	1/1	0.98	0.04	24,24,24,24	0
2	PGO	D	301	5/5	0.99	0.04	26,26,27,28	0
2	PGO	D	302	5/5	0.99	0.03	21,22,24,24	0
2	PGO	B	301	5/5	0.99	0.04	25,25,27,27	0
2	PGO	G	301	5/5	0.99	0.06	26,28,30,30	0
2	PGO	I	301	5/5	0.99	0.04	28,28,29,29	0
3	CL	A	304	1/1	0.99	0.03	31,31,31,31	0
3	CL	B	303	1/1	0.99	0.03	29,29,29,29	0
3	CL	B	304	1/1	0.99	0.03	27,27,27,27	0
3	CL	F	301	1/1	0.99	0.02	29,29,29,29	0
3	CL	F	302	1/1	0.99	0.02	30,30,30,30	0
3	CL	F	303	1/1	0.99	0.03	31,31,31,31	0
2	PGO	B	302	5/5	0.99	0.03	19,20,21,24	0
3	CL	H	301	1/1	0.99	0.02	36,36,36,36	0
3	CL	I	302	1/1	0.99	0.03	34,34,34,34	0
4	K	A	305	1/1	0.99	0.03	44,44,44,44	0
4	K	B	305	1/1	0.99	0.04	44,44,44,44	0
4	K	C	304	1/1	0.99	0.03	29,29,29,29	0
4	K	E	302	1/1	0.99	0.03	47,47,47,47	0
4	K	G	306	1/1	0.99	0.04	34,34,34,34	0
5	PGR	C	301	5/5	0.99	0.04	34,36,39,39	0
6	SO4	G	303	5/5	0.99	0.03	34,34,35,36	0
3	CL	I	303	1/1	1.00	0.01	21,21,21,21	0
3	CL	C	303	1/1	1.00	0.03	35,35,35,35	0
3	CL	E	301	1/1	1.00	0.04	27,27,27,27	0
3	CL	G	304	1/1	1.00	0.02	27,27,27,27	0
4	K	D	305	1/1	1.00	0.02	32,32,32,32	0
3	CL	G	305	1/1	1.00	0.02	30,30,30,30	0
3	CL	A	303	1/1	1.00	0.04	39,39,39,39	0
3	CL	H	302	1/1	1.00	0.02	24,24,24,24	0
6	SO4	D	304	5/5	1.00	0.02	29,30,32,35	0
6	SO4	G	302	5/5	1.00	0.04	26,27,27,28	0
3	CL	C	302	1/1	1.00	0.03	31,31,31,31	0

6.5 Other polymers

There are no such residues in this entry.