



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 09:45 AM UTC

PDB ID : 4I1Y / pdb_00004i1y
Title : The structure of Cysteine synthase from Mycobacterium ulcerans Agy99
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)
Deposited on : 2012-11-21
Resolution : 2.60 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

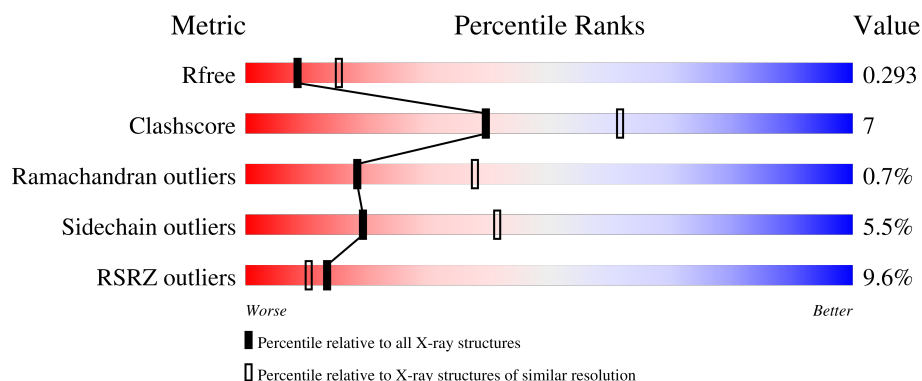
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	314	4% 71% 11% • 16%
1	B	314	9% 69% 12% • 16%
1	C	314	11% 73% 11% • 15%
1	D	314	8% 72% 12% •• 14%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7518 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 Cysteine synthase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	2	0
			1856	1156	325	365	10			
1	B	264	Total	C	N	O	S	0	0	0
			1846	1152	326	359	9			
1	C	268	Total	C	N	O	S	0	1	0
			1896	1191	326	370	9			
1	D	270	Total	C	N	O	S	0	0	0
			1863	1174	321	360	8			

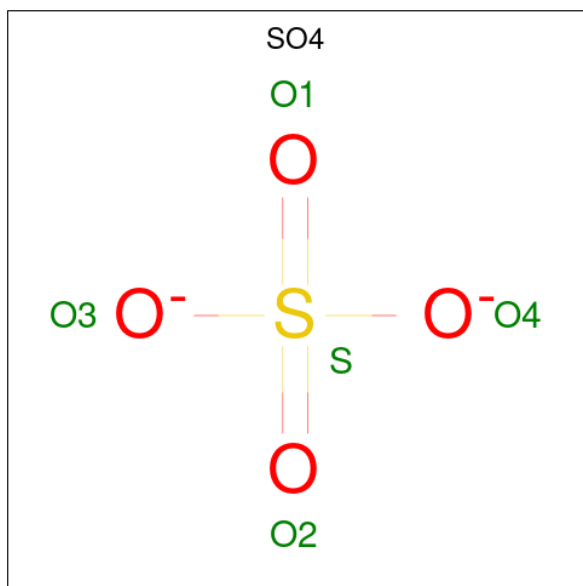
There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	GLY	-	expression tag	UNP A0PNS1
A	-2	PRO	-	expression tag	UNP A0PNS1
A	-1	GLY	-	expression tag	UNP A0PNS1
A	0	SER	-	expression tag	UNP A0PNS1
B	-3	GLY	-	expression tag	UNP A0PNS1
B	-2	PRO	-	expression tag	UNP A0PNS1
B	-1	GLY	-	expression tag	UNP A0PNS1
B	0	SER	-	expression tag	UNP A0PNS1
C	-3	GLY	-	expression tag	UNP A0PNS1
C	-2	PRO	-	expression tag	UNP A0PNS1
C	-1	GLY	-	expression tag	UNP A0PNS1
C	0	SER	-	expression tag	UNP A0PNS1
D	-3	GLY	-	expression tag	UNP A0PNS1
D	-2	PRO	-	expression tag	UNP A0PNS1
D	-1	GLY	-	expression tag	UNP A0PNS1
D	0	SER	-	expression tag	UNP A0PNS1

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Cl 1 1	0	0
2	B	1	Total Cl 1 1	0	0
2	C	1	Total Cl 1 1	0	0

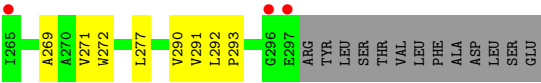
- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



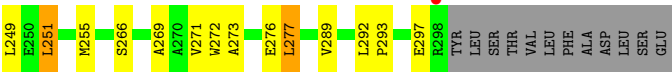
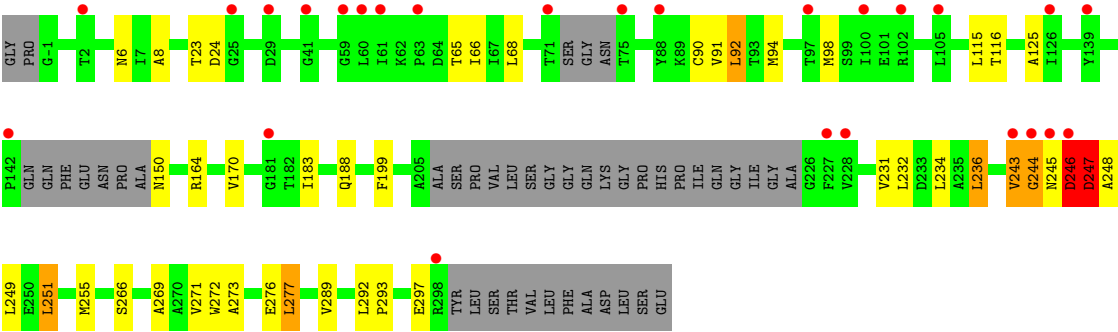
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	B	1	Total O S 5 4 1	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	8	Total O 8 8	0	0
4	B	17	Total O 17 17	0	0
4	C	10	Total O 10 10	0	0
4	D	9	Total O 9 9	0	0



● Molecule 1: Cysteine synthase



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	58.97Å 77.90Å 87.10Å 114.21° 104.86° 94.37°	Depositor
Resolution (Å)	48.71 – 2.60 48.71 – 2.60	Depositor EDS
% Data completeness (in resolution range)	97.5 (48.71-2.60) 97.9 (48.71-2.60)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.29 (at 2.58Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.261 , 0.293 0.256 , 0.293	Depositor DCC
R_{free} test set	2029 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	62.3	Xtriage
Anisotropy	0.251	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 53.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7518	wwPDB-VP
Average B, all atoms (Å ²)	82.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.12% 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: SO4, CL

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	0.53	0/1884	0.87	5/2566 (0.2%)
1	B	0.55	1/1868 (0.1%)	0.88	3/2544 (0.1%)
1	C	0.54	0/1925	0.87	4/2623 (0.2%)
1	D	0.61	1/1888 (0.1%)	0.85	2/2580 (0.1%)
All	All	0.56	2/7565 (0.0%)	0.87	14/10313 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	2
1	C	0	1
1	D	0	2
All	All	0	6

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	243	VAL	CA-C	-9.79	1.43	1.53
1	B	0	SER	C-N	-7.10	1.24	1.33

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	246	ASP	CB-CA-C	7.22	116.98	109.83
1	A	246	ASP	CB-CA-C	7.11	116.87	109.83
1	B	3	ILE	N-CA-C	6.15	116.94	107.78

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	246	ASP	CB-CA-C	5.98	115.75	109.83
1	D	246	ASP	O-C-N	-5.82	114.85	122.59
1	B	247	ASP	N-CA-C	-5.66	105.20	111.71
1	A	3	ILE	N-CA-C	5.54	116.54	107.24
1	C	247	ASP	N-CA-C	-5.53	105.35	111.71
1	C	150	ASN	CA-C-N	-5.49	114.27	119.76
1	C	150	ASN	C-N-CA	-5.49	114.27	119.76
1	A	242	THR	N-CA-C	-5.33	101.09	109.25
1	D	247	ASP	N-CA-C	5.17	121.82	110.80
1	A	246	ASP	N-CA-CB	-5.14	108.96	114.10
1	A	247	ASP	N-CA-C	-5.00	105.96	111.71

There are no chirality outliers.

All (6) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	246	ASP	Peptide
1	B	150	ASN	Peptide
1	B	246	ASP	Peptide
1	C	246	ASP	Peptide
1	D	150	ASN	Peptide
1	D	246	ASP	Peptide

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1856	0	1784	21	0
1	B	1846	0	1764	30	0
1	C	1896	0	1846	23	0
1	D	1863	0	1774	25	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
4	A	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	17	0	0	0	0
4	C	10	0	0	0	0
4	D	9	0	0	0	0
All	All	7518	0	7168	98	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 7.

All (98) 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:C:94:MET:HE1	1:C:113:LEU:HD22	1.67	0.75
1:A:94:MET:HE1	1:A:113:LEU:HD22	1.68	0.75
1:B:187:ALA:HA	1:B:199:PHE:CZ	2.22	0.75
1:D:246:ASP:HA	1:D:249:LEU:H	1.51	0.75
1:D:244:GLY:O	1:D:247:ASP:HB2	1.87	0.74
1:A:142:PRO:O	1:A:143:GLN:HB3	1.91	0.70
1:B:94:MET:HE1	1:B:113:LEU:HD22	1.73	0.69
1:A:243:VAL:O	1:A:245:ASN:N	2.28	0.65
1:C:188:GLN:HG3	1:C:236:LEU:HD11	1.79	0.63
1:D:243:VAL:O	1:D:245:ASN:N	2.33	0.62
1:D:243:VAL:CB	1:D:269:ALA:HB1	2.30	0.61
1:B:116:THR:HG21	1:B:125:ALA:HA	1.84	0.60
1:C:291:VAL:C	1:C:292:LEU:HD23	2.27	0.60
1:C:234:LEU:C	1:C:234:LEU:HD23	2.26	0.60
1:C:243:VAL:O	1:C:245:ASN:N	2.35	0.60
1:B:243:VAL:O	1:B:245:ASN:N	2.35	0.58
1:C:234:LEU:HD23	1:C:235:ALA:N	2.18	0.58
1:B:275:ARG:HH11	1:B:275:ARG:HG2	1.69	0.57
1:D:266:SER:N	1:D:297:GLU:OE1	2.36	0.57
1:D:94:MET:HE2	1:D:98:MET:CB	2.36	0.56
1:C:291:VAL:O	1:C:292:LEU:HD23	2.05	0.56
1:C:292:LEU:HA	1:C:293:PRO:C	2.32	0.55
1:B:261:LEU:HD23	1:C:7:ILE:HD12	1.89	0.54
1:D:251:LEU:HD23	1:D:272:TRP:CA	2.37	0.54
1:B:187:ALA:HA	1:B:199:PHE:CE2	2.43	0.54
1:A:69:GLU:OE2	1:A:142:PRO:HD2	2.08	0.54
1:C:116:THR:HG23	1:C:117:PRO:HD2	1.91	0.53
1:B:150:ASN:CB	1:B:151:PRO:CD	2.87	0.53
1:A:235:ALA:HB3	1:A:236:LEU:HD12	1.91	0.53
1:A:236:LEU:HD12	1:A:236:LEU:N	2.25	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:174:VAL:HB	1:B:290:VAL:HG12	1.90	0.52
1:A:116:THR:HG21	1:A:125:ALA:HA	1.93	0.51
1:B:246:ASP:HA	1:B:249:LEU:H	1.76	0.51
1:B:47:ILE:HG22	1:B:51:MET:HE2	1.94	0.50
1:A:292:LEU:HA	1:A:293:PRO:C	2.36	0.50
1:B:275:ARG:HH11	1:B:275:ARG:CG	2.24	0.50
1:C:251:LEU:HD23	1:C:272:TRP:CA	2.42	0.49
1:C:245:ASN:C	1:C:246:ASP:OD1	2.55	0.49
1:B:245:ASN:C	1:B:246:ASP:OD1	2.55	0.48
1:C:55:ALA:HB3	1:C:61:ILE:HD12	1.95	0.48
1:C:90:CYS:SG	1:C:92:LEU:HD13	2.53	0.48
1:D:6:ASN:OD1	1:D:8:ALA:HB3	2.13	0.48
1:D:170:VAL:HG23	1:D:289:VAL:HG23	1.95	0.48
1:A:68:LEU:O	1:A:141:ILE:HA	2.14	0.48
1:A:251:LEU:HD23	1:A:272:TRP:CA	2.43	0.48
1:D:116:THR:HG21	1:D:125:ALA:HA	1.96	0.48
1:D:65:THR:HG22	1:D:66:ILE:N	2.29	0.47
1:B:233:ASP:C	1:B:235:ALA:N	2.71	0.47
1:A:245:ASN:C	1:A:246:ASP:OD1	2.57	0.47
1:B:266:SER:CB	1:B:297:GLU:HG2	2.45	0.47
1:B:244:GLY:O	1:B:245:ASN:C	2.57	0.46
1:C:244:GLY:O	1:C:245:ASN:C	2.59	0.46
1:B:243:VAL:CB	1:B:269:ALA:HB1	2.46	0.46
1:B:251:LEU:CD2	1:B:272:TRP:HA	2.45	0.46
1:D:255:MET:SD	1:D:271:VAL:HG21	2.56	0.46
1:A:98:MET:HE2	1:A:98:MET:HB3	1.86	0.45
1:C:243:VAL:CB	1:C:269:ALA:HB1	2.47	0.45
1:D:183:ILE:HG22	1:D:199:PHE:CZ	2.52	0.45
1:D:292:LEU:HA	1:D:293:PRO:C	2.42	0.45
1:A:246:ASP:HA	1:A:249:LEU:H	1.82	0.44
1:B:150:ASN:CB	1:B:151:PRO:HD3	2.47	0.44
1:C:245:ASN:O	1:C:246:ASP:HB2	2.17	0.44
1:C:246:ASP:HA	1:C:249:LEU:H	1.82	0.44
1:D:243:VAL:O	1:D:244:GLY:C	2.60	0.44
1:D:251:LEU:CD2	1:D:272:TRP:HA	2.47	0.44
1:A:77:ILE:HD13	1:A:105:LEU:HD23	2.00	0.44
1:B:251:LEU:HD23	1:B:272:TRP:N	2.33	0.44
1:D:251:LEU:HD23	1:D:272:TRP:HA	2.00	0.44
1:A:23:THR:O	1:A:24:ASP:C	2.60	0.44
1:A:244:GLY:O	1:A:245:ASN:C	2.60	0.44
1:C:116:THR:HG22	1:C:117:PRO:O	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:267:SER:HB3	1:A:292:LEU:HD22	2.00	0.43
1:B:251:LEU:HD12	1:B:251:LEU:HA	1.81	0.43
1:A:244:GLY:O	1:A:247:ASP:HB2	2.19	0.43
1:B:23:THR:O	1:B:24:ASP:C	2.61	0.42
1:B:245:ASN:O	1:B:246:ASP:HB2	2.18	0.42
1:D:68:LEU:HA	1:D:91:VAL:O	2.18	0.42
1:B:187:ALA:HB2	1:B:199:PHE:CD2	2.54	0.42
1:A:245:ASN:O	1:A:246:ASP:HB2	2.20	0.42
1:A:255:MET:SD	1:A:271:VAL:HG21	2.60	0.42
1:D:245:ASN:O	1:D:246:ASP:CB	2.66	0.42
1:C:23:THR:O	1:C:24:ASP:C	2.63	0.42
1:C:251:LEU:CD2	1:C:272:TRP:HA	2.50	0.41
1:D:188:GLN:HG3	1:D:236:LEU:HD11	2.03	0.41
1:D:273:ALA:O	1:D:277:LEU:HD22	2.19	0.41
1:B:90:CYS:SG	1:B:92:LEU:HD13	2.61	0.41
1:B:68:LEU:O	1:B:141:ILE:HA	2.20	0.41
1:C:255:MET:SD	1:C:271:VAL:HG21	2.60	0.41
1:D:23:THR:O	1:D:24:ASP:C	2.63	0.41
1:D:246:ASP:H	1:D:248:ALA:H	1.69	0.41
1:B:77:ILE:HD13	1:B:105:LEU:CD2	2.50	0.41
1:A:170:VAL:HG23	1:A:289:VAL:HG23	2.03	0.41
1:C:68:LEU:O	1:C:141:ILE:HA	2.21	0.41
1:D:90:CYS:SG	1:D:92:LEU:HD13	2.62	0.40
1:D:65:THR:CG2	1:D:66:ILE:N	2.84	0.40
1:B:173:PHE:CD2	1:B:199:PHE:CE2	3.10	0.40
1:B:292:LEU:HA	1:B:293:PRO:C	2.46	0.40
1:B:71:THR:HG23	1:B:98:MET:HE1	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	259/314 (82%)	253 (98%)	5 (2%)	1 (0%)	30	51
1	B	256/314 (82%)	250 (98%)	5 (2%)	1 (0%)	30	51
1	C	263/314 (84%)	255 (97%)	6 (2%)	2 (1%)	16	34
1	D	262/314 (83%)	255 (97%)	4 (2%)	3 (1%)	11	25
All	All	1040/1256 (83%)	1013 (97%)	20 (2%)	7 (1%)	18	38

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	247	ASP
1	A	244	GLY
1	B	244	GLY
1	C	244	GLY
1	D	244	GLY
1	C	181	GLY
1	D	246	ASP

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	173/237 (73%)	164 (95%)	9 (5%)	21	44
1	B	168/237 (71%)	160 (95%)	8 (5%)	23	47
1	C	181/237 (76%)	170 (94%)	11 (6%)	17	37
1	D	168/237 (71%)	158 (94%)	10 (6%)	17	37
All	All	690/948 (73%)	652 (94%)	38 (6%)	19	41

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	75	THR
1	A	92	LEU
1	A	98	MET
1	A	130	GLU

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Mol	Chain	Res	Type
1	A	135	THR
1	A	143	GLN
1	A	251	LEU
1	A	276	GLU
1	A	277	LEU
1	B	75	THR
1	B	92	LEU
1	B	96	ASP
1	B	98	MET
1	B	275	ARG
1	B	276	GLU
1	B	277	LEU
1	B	297	GLU
1	C	75	THR
1	C	92	LEU
1	C	97	THR
1	C	98	MET
1	C	100	ILE
1	C	135	THR
1	C	234	LEU
1	C	236	LEU
1	C	251	LEU
1	C	277	LEU
1	C	290	VAL
1	D	92	LEU
1	D	115	LEU
1	D	164	ARG
1	D	231	VAL
1	D	232	LEU
1	D	234	LEU
1	D	236	LEU
1	D	251	LEU
1	D	276	GLU
1	D	277	LEU

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

Mol	Chain	Res	Type
1	A	143	GLN
1	A	198	GLN
1	C	198	GLN
1	D	198	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](#)

Of 5 ligands modelled in this entry, 3 are monoatomic - leaving 2 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	SO4	A	402	-	4,4,4	0.40	0	6,6,6	0.06	0
3	SO4	B	402	-	4,4,4	0.37	0	6,6,6	0.08	0

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 ⓘ

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	263/314 (83%)	0.54	13 (4%) 35 29	39, 71, 96, 256	2 (0%)
1	B	264/314 (84%)	0.76	27 (10%) 12 9	58, 77, 112, 294	0
1	C	268/314 (85%)	1.17	36 (13%) 7 5	51, 75, 179, 261	1 (0%)
1	D	270/314 (85%)	0.77	26 (9%) 13 10	61, 72, 131, 173	0
All	All	1065/1256 (84%)	0.81	102 (9%) 13 10	39, 74, 120, 294	3 (0%)

All (102) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	-2	PRO	10.4
1	C	156	VAL	9.2
1	A	297	GLU	8.7
1	C	153	VAL	8.6
1	C	154	HIS	8.1
1	C	155	ALA	7.8
1	C	158	THR	7.4
1	C	157	THR	7.3
1	C	168	GLY	7.2
1	C	0	SER	7.1
1	C	163	TRP	7.1
1	C	160	GLU	6.9
1	C	151	PRO	6.4
1	C	159	ALA	6.3
1	C	164	ARG	6.3
1	C	-1	GLY	6.2
1	C	161	GLU	6.2
1	C	162	VAL	6.1
1	C	165	ASP	6.0
1	C	152	ALA	6.0
1	C	150	ASN	5.9

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Mol	Chain	Res	Type	RSRZ
1	C	169	LYS	5.8
1	B	297	GLU	5.7
1	C	166	THR	5.6
1	D	75	THR	5.0
1	A	296	GLY	4.8
1	D	245	ASN	4.1
1	C	167	ASP	3.7
1	B	75	THR	3.7
1	D	97	THR	3.6
1	D	61	ILE	3.6
1	B	150	ASN	3.5
1	B	180	GLY	3.5
1	D	105	LEU	3.5
1	D	102	ARG	3.1
1	B	63	PRO	3.1
1	C	297	GLU	3.0
1	A	294	ASP	3.0
1	A	2	THR	3.0
1	D	100	ILE	3.0
1	C	75	THR	3.0
1	A	233	ASP	2.9
1	D	63	PRO	2.9
1	C	296	GLY	2.9
1	A	204	PRO	2.9
1	A	295	PHE	2.9
1	A	180	GLY	2.8
1	B	54	ALA	2.8
1	D	244	GLY	2.7
1	B	71	THR	2.7
1	B	1	MET	2.7
1	B	135	THR	2.7
1	D	71	THR	2.7
1	B	56	GLU	2.6
1	D	59	GLY	2.6
1	B	51	MET	2.6
1	B	53	ASP	2.6
1	B	294	ASP	2.6
1	A	243	VAL	2.6
1	C	245	ASN	2.6
1	A	242	THR	2.6
1	D	246	ASP	2.6
1	B	295	PHE	2.6

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Mol	Chain	Res	Type	RSRZ
1	B	61	ILE	2.5
1	D	298	ARG	2.5
1	B	74	ASN	2.5
1	B	246	ASP	2.5
1	B	88	TYR	2.5
1	C	180	GLY	2.5
1	C	244	GLY	2.5
1	D	2	THR	2.5
1	D	139	TYR	2.5
1	D	25	GLY	2.4
1	C	242	THR	2.4
1	B	81	MET	2.3
1	B	296	GLY	2.3
1	D	228	VAL	2.3
1	B	164	ARG	2.3
1	C	243	VAL	2.3
1	A	25	GLY	2.3
1	C	97	THR	2.2
1	B	76	GLY	2.2
1	D	181	GLY	2.2
1	C	265	ILE	2.2
1	D	227	PHE	2.2
1	D	88	TYR	2.2
1	C	87	GLY	2.2
1	A	246	ASP	2.2
1	D	29	ASP	2.2
1	A	177	VAL	2.2
1	C	199	PHE	2.2
1	D	142	PRO	2.1
1	B	257	THR	2.1
1	D	41	GLY	2.1
1	B	67	ILE	2.1
1	D	126	ILE	2.1
1	B	205	ALA	2.1
1	B	68	LEU	2.1
1	B	52	ILE	2.1
1	C	230	PRO	2.0
1	D	60	LEU	2.0
1	D	243	VAL	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 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	CL	A	401	1/1	-0.15	0.35	292,292,292,292	0
2	CL	C	401	1/1	-0.15	0.28	289,289,289,289	0
3	SO4	A	402	5/5	0.31	0.26	256,256,256,256	0
3	SO4	B	402	5/5	0.39	0.23	256,256,256,256	0
2	CL	B	401	1/1	0.63	0.38	304,304,304,304	0

6.5 Other polymers [i](#)

There are no such residues in this entry.