



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 14, 2026 – 02:12 PM UTC

PDB ID : 6OVT / pdb_00006ovt
Title : Crystal Structure of IlvD from Mycobacterium tuberculosis
Authors : Almo, S.C.; Grove, T.L.; Bonanno, J.B.; Baker, E.N.; Bashiri, G.
Deposited on : 2019-05-08
Resolution : 1.88 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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

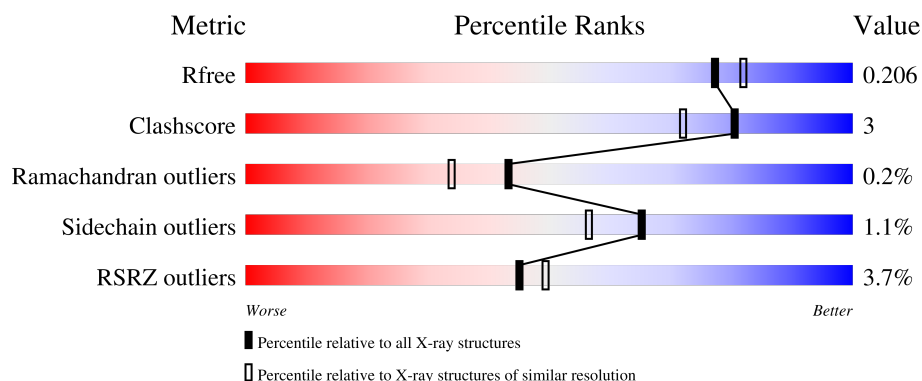
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.88 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	1220 (1.88-1.88)
Clashscore	190562	1234 (1.88-1.88)
Ramachandran outliers	187476	1222 (1.88-1.88)
Sidechain outliers	187428	1222 (1.88-1.88)
RSRZ outliers	180081	1220 (1.88-1.88)

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	600	
1	B	600	
1	C	600	
1	D	600	

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 17653 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 Dihydroxy-acid dehydratase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	562	Total	C	N	O	S	0	7	0
			4097	2551	727	794	25			
1	B	555	Total	C	N	O	S	0	3	0
			4041	2516	722	777	26			
1	C	558	Total	C	N	O	S	0	10	0
			4098	2551	731	791	25			
1	D	560	Total	C	N	O	S	0	10	0
			4109	2557	732	794	26			

There are 100 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-24	MET	-	initiating methionine	UNP A0A0E8UWV6
A	-23	SER	-	expression tag	UNP A0A0E8UWV6
A	-22	TYR	-	expression tag	UNP A0A0E8UWV6
A	-21	TYR	-	expression tag	UNP A0A0E8UWV6
A	-20	HIS	-	expression tag	UNP A0A0E8UWV6
A	-19	HIS	-	expression tag	UNP A0A0E8UWV6
A	-18	HIS	-	expression tag	UNP A0A0E8UWV6
A	-17	HIS	-	expression tag	UNP A0A0E8UWV6
A	-16	HIS	-	expression tag	UNP A0A0E8UWV6
A	-15	HIS	-	expression tag	UNP A0A0E8UWV6
A	-14	ASP	-	expression tag	UNP A0A0E8UWV6
A	-13	TYR	-	expression tag	UNP A0A0E8UWV6
A	-12	ASP	-	expression tag	UNP A0A0E8UWV6
A	-11	ILE	-	expression tag	UNP A0A0E8UWV6
A	-10	PRO	-	expression tag	UNP A0A0E8UWV6
A	-9	THR	-	expression tag	UNP A0A0E8UWV6
A	-8	THR	-	expression tag	UNP A0A0E8UWV6
A	-7	GLU	-	expression tag	UNP A0A0E8UWV6
A	-6	ASN	-	expression tag	UNP A0A0E8UWV6
A	-5	LEU	-	expression tag	UNP A0A0E8UWV6
A	-4	TYR	-	expression tag	UNP A0A0E8UWV6

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
A	-3	PHE	-	expression tag	UNP A0A0E8UWV6
A	-2	GLN	-	expression tag	UNP A0A0E8UWV6
A	-1	GLY	-	expression tag	UNP A0A0E8UWV6
A	0	ALA	-	expression tag	UNP A0A0E8UWV6
B	-24	MET	-	initiating methionine	UNP A0A0E8UWV6
B	-23	SER	-	expression tag	UNP A0A0E8UWV6
B	-22	TYR	-	expression tag	UNP A0A0E8UWV6
B	-21	TYR	-	expression tag	UNP A0A0E8UWV6
B	-20	HIS	-	expression tag	UNP A0A0E8UWV6
B	-19	HIS	-	expression tag	UNP A0A0E8UWV6
B	-18	HIS	-	expression tag	UNP A0A0E8UWV6
B	-17	HIS	-	expression tag	UNP A0A0E8UWV6
B	-16	HIS	-	expression tag	UNP A0A0E8UWV6
B	-15	HIS	-	expression tag	UNP A0A0E8UWV6
B	-14	ASP	-	expression tag	UNP A0A0E8UWV6
B	-13	TYR	-	expression tag	UNP A0A0E8UWV6
B	-12	ASP	-	expression tag	UNP A0A0E8UWV6
B	-11	ILE	-	expression tag	UNP A0A0E8UWV6
B	-10	PRO	-	expression tag	UNP A0A0E8UWV6
B	-9	THR	-	expression tag	UNP A0A0E8UWV6
B	-8	THR	-	expression tag	UNP A0A0E8UWV6
B	-7	GLU	-	expression tag	UNP A0A0E8UWV6
B	-6	ASN	-	expression tag	UNP A0A0E8UWV6
B	-5	LEU	-	expression tag	UNP A0A0E8UWV6
B	-4	TYR	-	expression tag	UNP A0A0E8UWV6
B	-3	PHE	-	expression tag	UNP A0A0E8UWV6
B	-2	GLN	-	expression tag	UNP A0A0E8UWV6
B	-1	GLY	-	expression tag	UNP A0A0E8UWV6
B	0	ALA	-	expression tag	UNP A0A0E8UWV6
C	-24	MET	-	initiating methionine	UNP A0A0E8UWV6
C	-23	SER	-	expression tag	UNP A0A0E8UWV6
C	-22	TYR	-	expression tag	UNP A0A0E8UWV6
C	-21	TYR	-	expression tag	UNP A0A0E8UWV6
C	-20	HIS	-	expression tag	UNP A0A0E8UWV6
C	-19	HIS	-	expression tag	UNP A0A0E8UWV6
C	-18	HIS	-	expression tag	UNP A0A0E8UWV6
C	-17	HIS	-	expression tag	UNP A0A0E8UWV6
C	-16	HIS	-	expression tag	UNP A0A0E8UWV6
C	-15	HIS	-	expression tag	UNP A0A0E8UWV6
C	-14	ASP	-	expression tag	UNP A0A0E8UWV6
C	-13	TYR	-	expression tag	UNP A0A0E8UWV6
C	-12	ASP	-	expression tag	UNP A0A0E8UWV6

Continued on next page...

Continued from previous page...

Chain	Residue	Modelled	Actual	Comment	Reference
C	-11	ILE	-	expression tag	UNP A0A0E8UWV6
C	-10	PRO	-	expression tag	UNP A0A0E8UWV6
C	-9	THR	-	expression tag	UNP A0A0E8UWV6
C	-8	THR	-	expression tag	UNP A0A0E8UWV6
C	-7	GLU	-	expression tag	UNP A0A0E8UWV6
C	-6	ASN	-	expression tag	UNP A0A0E8UWV6
C	-5	LEU	-	expression tag	UNP A0A0E8UWV6
C	-4	TYR	-	expression tag	UNP A0A0E8UWV6
C	-3	PHE	-	expression tag	UNP A0A0E8UWV6
C	-2	GLN	-	expression tag	UNP A0A0E8UWV6
C	-1	GLY	-	expression tag	UNP A0A0E8UWV6
C	0	ALA	-	expression tag	UNP A0A0E8UWV6
D	-24	MET	-	initiating methionine	UNP A0A0E8UWV6
D	-23	SER	-	expression tag	UNP A0A0E8UWV6
D	-22	TYR	-	expression tag	UNP A0A0E8UWV6
D	-21	TYR	-	expression tag	UNP A0A0E8UWV6
D	-20	HIS	-	expression tag	UNP A0A0E8UWV6
D	-19	HIS	-	expression tag	UNP A0A0E8UWV6
D	-18	HIS	-	expression tag	UNP A0A0E8UWV6
D	-17	HIS	-	expression tag	UNP A0A0E8UWV6
D	-16	HIS	-	expression tag	UNP A0A0E8UWV6
D	-15	HIS	-	expression tag	UNP A0A0E8UWV6
D	-14	ASP	-	expression tag	UNP A0A0E8UWV6
D	-13	TYR	-	expression tag	UNP A0A0E8UWV6
D	-12	ASP	-	expression tag	UNP A0A0E8UWV6
D	-11	ILE	-	expression tag	UNP A0A0E8UWV6
D	-10	PRO	-	expression tag	UNP A0A0E8UWV6
D	-9	THR	-	expression tag	UNP A0A0E8UWV6
D	-8	THR	-	expression tag	UNP A0A0E8UWV6
D	-7	GLU	-	expression tag	UNP A0A0E8UWV6
D	-6	ASN	-	expression tag	UNP A0A0E8UWV6
D	-5	LEU	-	expression tag	UNP A0A0E8UWV6
D	-4	TYR	-	expression tag	UNP A0A0E8UWV6
D	-3	PHE	-	expression tag	UNP A0A0E8UWV6
D	-2	GLN	-	expression tag	UNP A0A0E8UWV6
D	-1	GLY	-	expression tag	UNP A0A0E8UWV6
D	0	ALA	-	expression tag	UNP A0A0E8UWV6

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

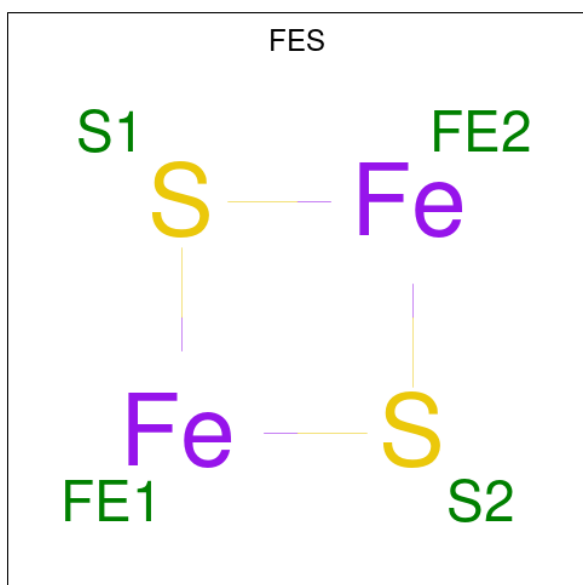
Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
-----	-------	----------	-------	---------	---------

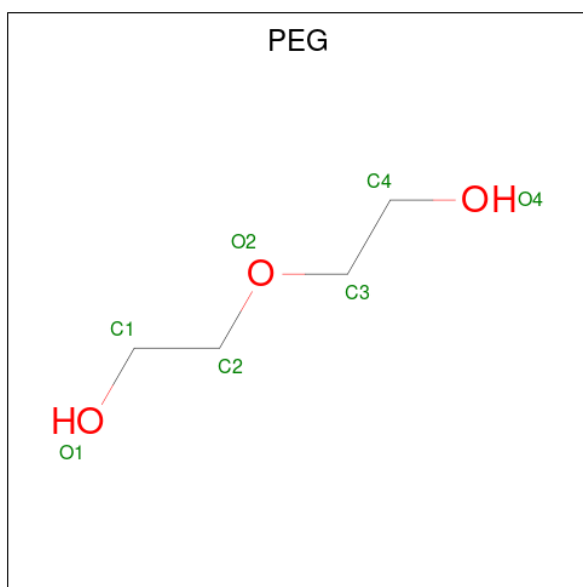
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	3	Total Mg 3 3	0	0
2	B	3	Total Mg 3 3	0	0
2	C	3	Total Mg 3 3	0	0
2	D	3	Total Mg 3 3	0	0

- Molecule 3 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Fe S 4 2 2	0	0
3	B	1	Total Fe S 4 2 2	0	0
3	C	1	Total Fe S 4 2 2	0	0
3	D	1	Total Fe S 4 2 2	0	0

- Molecule 4 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: $\text{C}_4\text{H}_{10}\text{O}_3$).



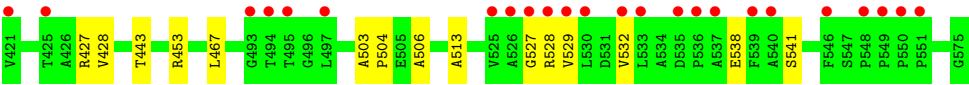
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			7	4	3		
4	B	1	Total	C	O	0	0
			7	4	3		
4	C	1	Total	C	O	0	0
			7	4	3		

- Molecule 5 is SODIUM ION (CCD ID: NA) (formula: Na).

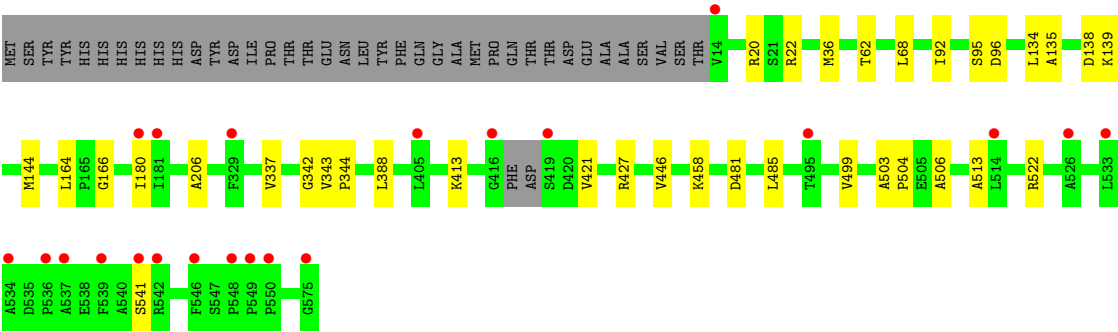
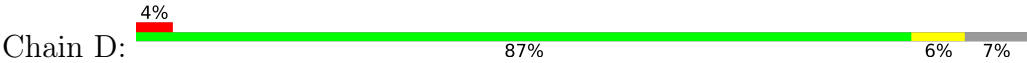
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	D	1	Total	Na	0	0
			1	1		

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	382	Total	O	0	0
			382	382		
6	B	256	Total	O	0	0
			256	256		
6	C	306	Total	O	0	0
			306	306		
6	D	314	Total	O	0	0
			314	314		



● Molecule 1: Dihydroxy-acid dehydratase



4 Data and refinement statistics

Property	Value	Source
Space group	P 61	Depositor
Cell constants a, b, c, α , β , γ	88.41Å 88.41Å 483.31Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	80.55 – 1.88 80.55 – 1.88	Depositor EDS
% Data completeness (in resolution range)	100.0 (80.55-1.88) 100.0 (80.55-1.88)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.95 (at 1.88Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.160 , 0.201 0.168 , 0.206	Depositor DCC
R_{free} test set	8430 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å ²)	26.7	Xtriage
Anisotropy	0.126	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 42.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.059 for h,-h-k,-l	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	17653	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.35% 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: PEG, NA, KCX, FES, MG

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	1.14	2/4163 (0.0%)	1.24	1/5643 (0.0%)
1	B	1.10	1/4092 (0.0%)	1.29	1/5541 (0.0%)
1	C	1.09	1/4165 (0.0%)	1.32	6/5641 (0.1%)
1	D	1.10	0/4176	1.28	1/5657 (0.0%)
All	All	1.10	4/16596 (0.0%)	1.28	9/22482 (0.0%)

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	2
1	D	0	1
All	All	0	3

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	69	ASP	C-O	5.23	1.30	1.24
1	A	53	GLY	C-O	5.13	1.29	1.24
1	C	296	HIS	CE1-NE2	5.10	1.37	1.32
1	A	50	PRO	C-O	-5.04	1.18	1.23

The worst 5 of 9 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	377	PRO	CB-CA-C	-5.93	105.81	111.39
1	B	179	THR	CA-CB-OG1	-5.89	100.77	109.60
1	C	157	PHE	CA-CB-CG	-5.77	108.03	113.80

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	171	SER	CA-C-N	5.71	128.72	120.38
1	C	171	SER	C-N-CA	5.71	128.72	120.38

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	214	CYS	Peptide
1	A	92	ILE	Peptide
1	D	92	ILE	Peptide

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	4097	0	4112	19	0
1	B	4041	0	4057	26	0
1	C	4098	0	4127	26	0
1	D	4109	0	4132	20	0
2	A	3	0	0	0	0
2	B	3	0	0	0	0
2	C	3	0	0	0	0
2	D	3	0	0	0	0
3	A	4	0	0	0	0
3	B	4	0	0	0	0
3	C	4	0	0	0	0
3	D	4	0	0	0	0
4	A	7	0	10	0	0
4	B	7	0	10	3	0
4	C	7	0	10	2	0
5	D	1	0	0	0	0
6	A	382	0	0	3	0
6	B	256	0	0	2	0
6	C	306	0	0	4	0
6	D	314	0	0	4	0
All	All	17653	0	16458	86	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 3.

The worst 5 of 86 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:427[A]:ARG:HE	1:C:443:THR:HG22	1.63	0.64
1:C:43:ASP:HB3	1:C:45:GLU:OE2	2.01	0.61
1:D:446:VAL:HG13	1:D:481:ASP:HB3	1.84	0.60
1:D:68:LEU:HD22	1:D:134:LEU:HB3	1.85	0.59
1:D:20[A]:ARG:NH2	6:D:702:HOH:O	2.16	0.58

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	566/600 (94%)	556 (98%)	9 (2%)	1 (0%)	43	34
1	B	551/600 (92%)	537 (98%)	13 (2%)	1 (0%)	43	34
1	C	563/600 (94%)	552 (98%)	10 (2%)	1 (0%)	43	34
1	D	565/600 (94%)	553 (98%)	11 (2%)	1 (0%)	43	34
All	All	2245/2400 (94%)	2198 (98%)	43 (2%)	4 (0%)	43	34

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	138	ASP
1	C	138	ASP
1	D	138	ASP
1	A	138	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	424/451 (94%)	419 (99%)	5 (1%)	63	53
1	B	415/451 (92%)	409 (99%)	6 (1%)	59	48
1	C	424/451 (94%)	419 (99%)	5 (1%)	63	53
1	D	425/451 (94%)	422 (99%)	3 (1%)	76	69
All	All	1688/1804 (94%)	1669 (99%)	19 (1%)	65	56

5 of 19 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	467	LEU
1	D	413	LYS
1	D	541	SER
1	D	388	LEU
1	B	392	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	402	HIS
1	A	543	GLN
1	C	390	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](#)

4 non-standard protein/DNA/RNA residues 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
1	KCX	A	139	1,2	10,11,12	1.49	1 (10%)	6,12,14	2.58	1 (16%)
1	KCX	C	139	1,2	10,11,12	0.54	0	6,12,14	0.77	0
1	KCX	D	139	1,2	10,11,12	2.37	1 (10%)	6,12,14	2.61	1 (16%)
1	KCX	B	139	1,2	10,11,12	0.49	0	6,12,14	0.90	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
1	KCX	A	139	1,2	-	0/9/10/12	-
1	KCX	C	139	1,2	-	0/9/10/12	-
1	KCX	D	139	1,2	-	0/9/10/12	-
1	KCX	B	139	1,2	-	0/9/10/12	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	139	KCX	OQ1-CX	7.09	1.34	1.21
1	A	139	KCX	OQ1-CX	4.29	1.29	1.21

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	139	KCX	OQ1-CX-NZ	-6.39	115.21	124.92
1	A	139	KCX	OQ1-CX-NZ	-6.28	115.38	124.92

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 20 ligands modelled in this entry, 13 are monoatomic - leaving 7 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	FES	C	601	1,6	0,4,4	-	-	-		
4	PEG	C	602	-	6,6,6	0.24	0	5,5,5	0.18	0
3	FES	D	603	1,6	0,4,4	-	-	-		
4	PEG	A	605	-	6,6,6	0.32	0	5,5,5	0.12	0
4	PEG	B	603	-	6,6,6	0.25	0	5,5,5	0.16	0
3	FES	A	602	1,6	0,4,4	-	-	-		
3	FES	B	602	1,6	0,4,4	-	-	-		

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	FES	C	601	1,6	-	-	0/1/1/1
4	PEG	C	602	-	-	1/4/4/4	-
3	FES	D	603	1,6	-	-	0/1/1/1
4	PEG	A	605	-	-	2/4/4/4	-
4	PEG	B	603	-	-	3/4/4/4	-
3	FES	A	602	1,6	-	-	0/1/1/1
3	FES	B	602	1,6	-	-	0/1/1/1

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 6 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	603	PEG	O1-C1-C2-O2
4	C	602	PEG	O2-C3-C4-O4
4	A	605	PEG	O1-C1-C2-O2
4	B	603	PEG	C4-C3-O2-C2
4	A	605	PEG	O2-C3-C4-O4

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	602	PEG	2	0
4	B	603	PEG	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	561/600 (93%)	-0.24	2 (0%) 88 92	10, 25, 45, 73	7 (1%)
1	B	554/600 (92%)	0.23	22 (3%) 42 45	12, 34, 62, 81	3 (0%)
1	C	557/600 (92%)	0.10	36 (6%) 25 27	11, 29, 68, 94	10 (1%)
1	D	559/600 (93%)	0.21	22 (3%) 43 47	11, 31, 62, 107	10 (1%)
All	All	2231/2400 (92%)	0.08	82 (3%) 45 49	10, 29, 60, 107	30 (1%)

The worst 5 of 82 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	415	ALA	5.3
1	D	14	VAL	4.5
1	B	15	ALA	4.2
1	C	184	PHE	4.2
1	D	575	GLY	4.2

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

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
1	KCX	C	139	12/13	0.92	0.10	22,23,30,32	0
1	KCX	B	139	12/13	0.94	0.09	27,30,34,35	0
1	KCX	D	139	12/13	0.94	0.08	24,27,34,35	0
1	KCX	A	139	12/13	0.96	0.06	18,19,23,23	0

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
5	NA	D	605	1/1	0.74	0.18	62,62,62,62	0
4	PEG	B	603	7/7	0.87	0.11	56,56,58,59	0
4	PEG	A	605	7/7	0.89	0.12	38,46,47,49	0
4	PEG	C	602	7/7	0.90	0.12	48,51,54,56	0
2	MG	D	604	1/1	0.93	0.06	49,49,49,49	0
2	MG	D	602	1/1	0.96	0.08	48,48,48,48	0
2	MG	A	603	1/1	0.96	0.04	50,50,50,50	0
2	MG	B	604	1/1	0.97	0.04	50,50,50,50	0
2	MG	D	601	1/1	0.97	0.09	40,40,40,40	0
2	MG	A	601	1/1	0.98	0.06	23,23,23,23	0
2	MG	C	603	1/1	0.98	0.12	32,32,32,32	0
2	MG	B	601	1/1	0.98	0.08	36,36,36,36	0
3	FES	B	602	4/4	0.99	0.04	28,28,30,31	0
3	FES	C	601	4/4	0.99	0.05	24,26,29,30	0
3	FES	D	603	4/4	0.99	0.04	23,25,25,27	0
2	MG	C	605	1/1	0.99	0.06	39,39,39,39	0
2	MG	B	605	1/1	0.99	0.05	48,48,48,48	0
2	MG	A	604	1/1	0.99	0.05	42,42,42,42	0
2	MG	C	604	1/1	0.99	0.04	41,41,41,41	0
3	FES	A	602	4/4	1.00	0.03	21,22,23,24	0

6.5 Other polymers ⓘ

There are no such residues in this entry.