



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:06 PM UTC

PDB ID : 5V2D / pdb_00005v2d
Title : Crystal structure of Pseudomonas brassicacearum lignostilbene dioxygenase
Authors : Loewen, P.C.; Loewen, M.
Deposited on : 2017-03-03
Resolution : 1.90 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
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

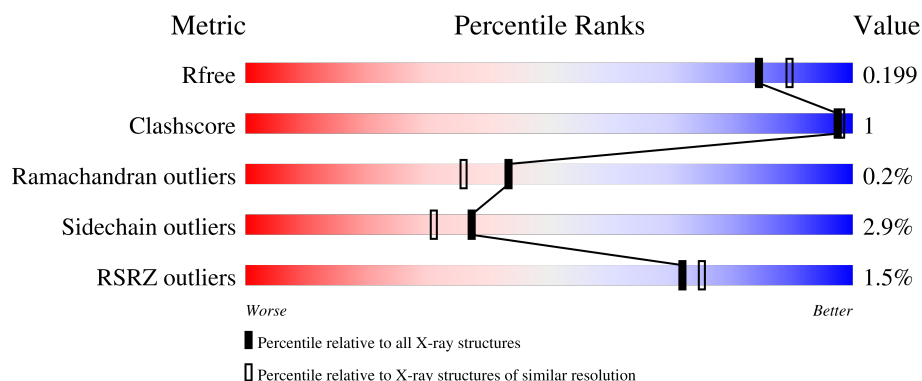
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	501	0% 90% 7% ..
1	B	501	0% 91% 5% .
1	C	501	3% 87% 8% ..
1	D	501	0% 90% 5% .

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 16428 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 Dioxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	488	Total	C	N	O	S	0	3	0
			3884	2495	662	710	17			
1	B	483	Total	C	N	O	S	0	4	0
			3850	2475	652	706	17			
1	C	480	Total	C	N	O	S	0	0	0
			3799	2443	642	698	16			
1	D	480	Total	C	N	O	S	0	2	0
			3815	2456	645	698	16			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP W8QAY8
A	-18	GLY	-	expression tag	UNP W8QAY8
A	-17	SER	-	expression tag	UNP W8QAY8
A	-16	SER	-	expression tag	UNP W8QAY8
A	-15	HIS	-	expression tag	UNP W8QAY8
A	-14	HIS	-	expression tag	UNP W8QAY8
A	-13	HIS	-	expression tag	UNP W8QAY8
A	-12	HIS	-	expression tag	UNP W8QAY8
A	-11	HIS	-	expression tag	UNP W8QAY8
A	-10	HIS	-	expression tag	UNP W8QAY8
A	-9	SER	-	expression tag	UNP W8QAY8
A	-8	SER	-	expression tag	UNP W8QAY8
A	-7	GLY	-	expression tag	UNP W8QAY8
A	-6	LEU	-	expression tag	UNP W8QAY8
A	-5	VAL	-	expression tag	UNP W8QAY8
A	-4	PRO	-	expression tag	UNP W8QAY8
A	-3	ARG	-	expression tag	UNP W8QAY8
A	-2	GLY	-	expression tag	UNP W8QAY8
A	-1	SER	-	expression tag	UNP W8QAY8
A	0	HIS	-	expression tag	UNP W8QAY8
B	-19	MET	-	expression tag	UNP W8QAY8

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-18	GLY	-	expression tag	UNP W8QAY8
B	-17	SER	-	expression tag	UNP W8QAY8
B	-16	SER	-	expression tag	UNP W8QAY8
B	-15	HIS	-	expression tag	UNP W8QAY8
B	-14	HIS	-	expression tag	UNP W8QAY8
B	-13	HIS	-	expression tag	UNP W8QAY8
B	-12	HIS	-	expression tag	UNP W8QAY8
B	-11	HIS	-	expression tag	UNP W8QAY8
B	-10	HIS	-	expression tag	UNP W8QAY8
B	-9	SER	-	expression tag	UNP W8QAY8
B	-8	SER	-	expression tag	UNP W8QAY8
B	-7	GLY	-	expression tag	UNP W8QAY8
B	-6	LEU	-	expression tag	UNP W8QAY8
B	-5	VAL	-	expression tag	UNP W8QAY8
B	-4	PRO	-	expression tag	UNP W8QAY8
B	-3	ARG	-	expression tag	UNP W8QAY8
B	-2	GLY	-	expression tag	UNP W8QAY8
B	-1	SER	-	expression tag	UNP W8QAY8
B	0	HIS	-	expression tag	UNP W8QAY8
C	-19	MET	-	expression tag	UNP W8QAY8
C	-18	GLY	-	expression tag	UNP W8QAY8
C	-17	SER	-	expression tag	UNP W8QAY8
C	-16	SER	-	expression tag	UNP W8QAY8
C	-15	HIS	-	expression tag	UNP W8QAY8
C	-14	HIS	-	expression tag	UNP W8QAY8
C	-13	HIS	-	expression tag	UNP W8QAY8
C	-12	HIS	-	expression tag	UNP W8QAY8
C	-11	HIS	-	expression tag	UNP W8QAY8
C	-10	HIS	-	expression tag	UNP W8QAY8
C	-9	SER	-	expression tag	UNP W8QAY8
C	-8	SER	-	expression tag	UNP W8QAY8
C	-7	GLY	-	expression tag	UNP W8QAY8
C	-6	LEU	-	expression tag	UNP W8QAY8
C	-5	VAL	-	expression tag	UNP W8QAY8
C	-4	PRO	-	expression tag	UNP W8QAY8
C	-3	ARG	-	expression tag	UNP W8QAY8
C	-2	GLY	-	expression tag	UNP W8QAY8
C	-1	SER	-	expression tag	UNP W8QAY8
C	0	HIS	-	expression tag	UNP W8QAY8
D	-19	MET	-	expression tag	UNP W8QAY8
D	-18	GLY	-	expression tag	UNP W8QAY8
D	-17	SER	-	expression tag	UNP W8QAY8

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-16	SER	-	expression tag	UNP W8QAY8
D	-15	HIS	-	expression tag	UNP W8QAY8
D	-14	HIS	-	expression tag	UNP W8QAY8
D	-13	HIS	-	expression tag	UNP W8QAY8
D	-12	HIS	-	expression tag	UNP W8QAY8
D	-11	HIS	-	expression tag	UNP W8QAY8
D	-10	HIS	-	expression tag	UNP W8QAY8
D	-9	SER	-	expression tag	UNP W8QAY8
D	-8	SER	-	expression tag	UNP W8QAY8
D	-7	GLY	-	expression tag	UNP W8QAY8
D	-6	LEU	-	expression tag	UNP W8QAY8
D	-5	VAL	-	expression tag	UNP W8QAY8
D	-4	PRO	-	expression tag	UNP W8QAY8
D	-3	ARG	-	expression tag	UNP W8QAY8
D	-2	GLY	-	expression tag	UNP W8QAY8
D	-1	SER	-	expression tag	UNP W8QAY8
D	0	HIS	-	expression tag	UNP W8QAY8

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

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

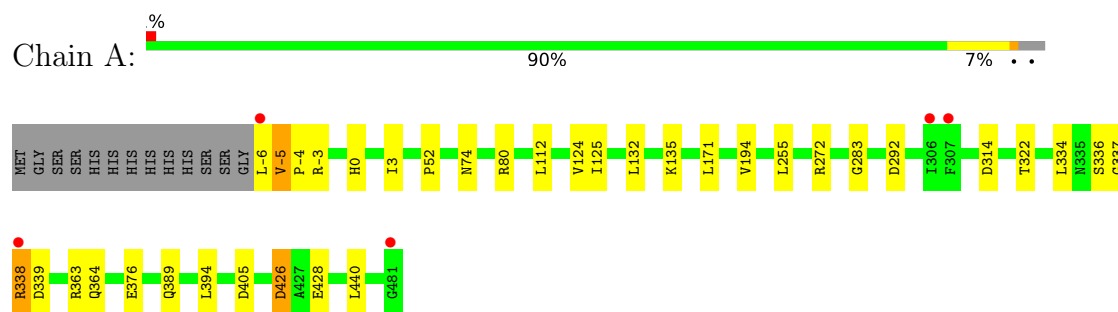
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	268	Total O 268 268	0	0
3	B	296	Total O 296 296	0	0
3	C	233	Total O 233 233	0	0
3	D	279	Total O 279 279	0	0

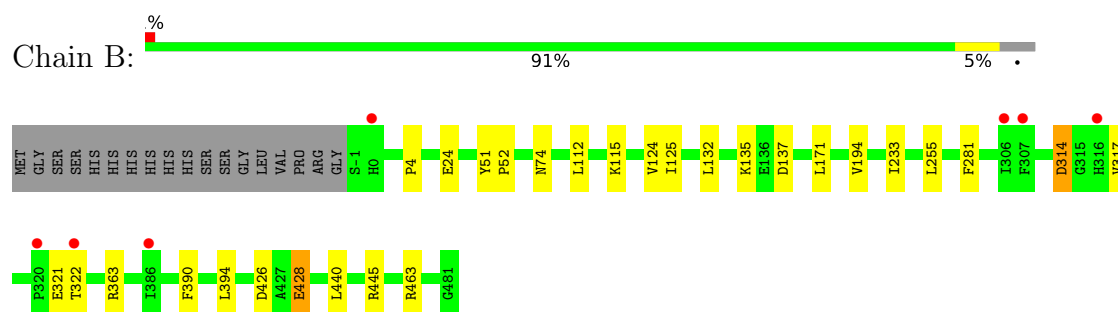
3 Residue-property plots [i](#)

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.

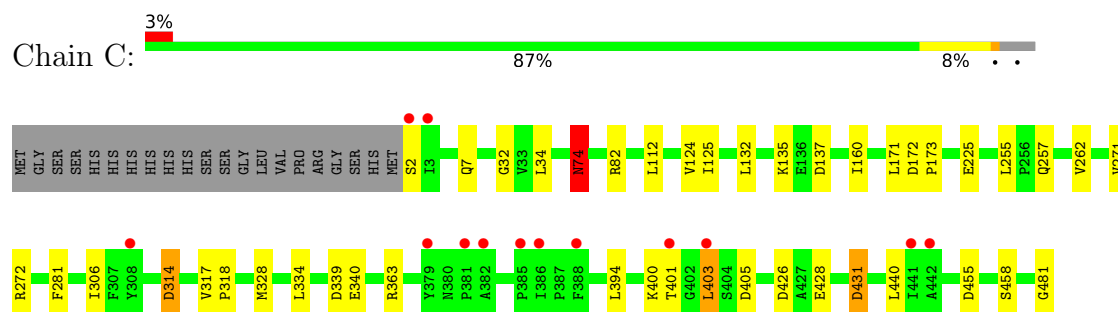
• Molecule 1: Dioxygenase



• Molecule 1: Dioxygenase

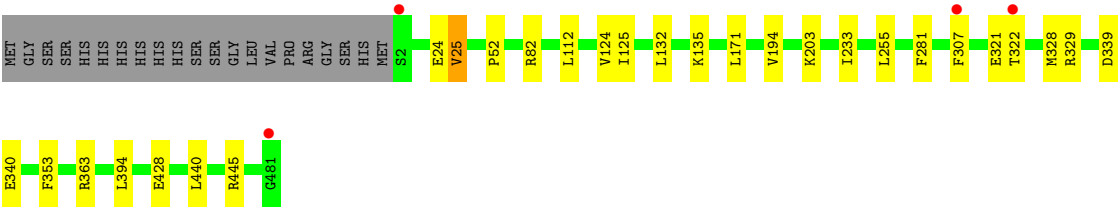


• Molecule 1: Dioxygenase



• Molecule 1: Dioxygenase





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	96.40Å 104.67Å 104.71Å 90.00° 94.79° 90.00°	Depositor
Resolution (Å)	48.03 – 1.90 48.03 – 1.90	Depositor EDS
% Data completeness (in resolution range)	98.8 (48.03-1.90) 98.8 (48.03-1.90)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.00 (at 1.90Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.161 , 0.193 0.171 , 0.199	Depositor DCC
R_{free} test set	8107 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	28.1	Xtriage
Anisotropy	0.366	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 34.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.97	EDS
Total number of atoms	16428	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.09% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: FE

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	1.35	12/4013 (0.3%)	1.15	11/5471 (0.2%)
1	B	1.32	6/3985 (0.2%)	1.13	9/5433 (0.2%)
1	C	1.37	11/3920 (0.3%)	1.12	12/5347 (0.2%)
1	D	1.31	5/3943 (0.1%)	1.11	8/5377 (0.1%)
All	All	1.34	34/15861 (0.2%)	1.13	40/21628 (0.2%)

All (34) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	-5	VAL	CA-CB	7.75	1.59	1.53
1	C	160	ILE	C-O	-6.87	1.16	1.24
1	C	281	PHE	C-O	6.64	1.31	1.23
1	A	405	ASP	CG-OD2	6.49	1.37	1.25
1	A	292	ASP	C-O	6.46	1.30	1.24
1	B	52	PRO	CA-C	6.42	1.58	1.52
1	B	51	TYR	N-CA	6.38	1.51	1.45
1	A	0	HIS	CA-C	6.23	1.61	1.52
1	C	271	VAL	C-O	-6.14	1.17	1.24
1	D	328	MET	SD-CE	-6.03	1.64	1.79
1	A	74	ASN	CB-CG	-5.97	1.37	1.52
1	A	-3	ARG	N-CA	5.97	1.53	1.46
1	C	172	ASP	CG-OD2	-5.90	1.14	1.25
1	C	340	GLU	C-O	-5.88	1.16	1.23
1	D	353	PHE	C-O	5.82	1.26	1.23
1	B	463	ARG	CZ-NH2	-5.74	1.25	1.33
1	A	283	GLY	N-CA	5.74	1.51	1.45
1	A	135	LYS	CG-CD	-5.65	1.35	1.52
1	A	0	HIS	C-O	-5.62	1.16	1.24
1	D	135	LYS	CG-CD	-5.48	1.36	1.52
1	C	262	VAL	N-CA	-5.48	1.40	1.46
1	B	4	PRO	CA-C	5.46	1.59	1.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	389	GLN	CA-C	5.44	1.60	1.53
1	C	318	PRO	CA-C	5.42	1.55	1.52
1	C	173	PRO	C-O	-5.41	1.17	1.24
1	B	4	PRO	C-O	-5.37	1.17	1.23
1	C	7	GLN	C-O	-5.22	1.17	1.24
1	D	52	PRO	CA-C	5.21	1.57	1.52
1	A	337	GLY	C-O	-5.19	1.18	1.24
1	C	32	GLY	C-O	-5.17	1.17	1.23
1	A	376	GLU	C-O	-5.09	1.17	1.24
1	D	329	ARG	CA-C	-5.09	1.46	1.52
1	C	431	ASP	C-O	-5.07	1.18	1.24
1	B	390	PHE	CA-C	5.03	1.59	1.52

All (40) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	25	VAL	CB-CA-C	9.22	122.98	110.99
1	C	314	ASP	N-CA-CB	-8.01	96.95	110.49
1	B	74	ASN	CB-CA-C	7.92	124.29	112.09
1	B	363	ARG	CG-CD-NE	7.91	129.40	112.00
1	A	339	ASP	N-CA-CB	-7.89	98.06	110.28
1	B	314	ASP	N-CA-CB	-7.55	97.74	110.49
1	A	338	ARG	CA-C-N	7.40	131.56	120.31
1	A	338	ARG	C-N-CA	7.40	131.56	120.31
1	B	322	THR	CB-CA-C	-7.30	97.12	110.63
1	D	322	THR	CB-CA-C	-7.27	97.18	110.63
1	C	281	PHE	N-CA-CB	-7.20	99.03	110.77
1	A	322	THR	CB-CA-C	-7.07	97.55	110.63
1	D	339	ASP	N-CA-CB	-6.95	100.31	110.53
1	D	25	VAL	CG1-CB-CG2	6.63	125.38	110.80
1	A	338	ARG	CA-C-O	6.54	127.51	120.38
1	D	194	VAL	CB-CA-C	-6.00	100.33	110.71
1	C	272	ARG	NE-CZ-NH2	5.99	124.59	119.20
1	B	194	VAL	CB-CA-C	-5.93	100.90	110.28
1	B	317	VAL	CA-C-N	-5.81	115.47	119.66
1	B	317	VAL	C-N-CA	-5.81	115.47	119.66
1	C	317	VAL	CA-C-N	-5.74	115.53	119.66
1	C	317	VAL	C-N-CA	-5.74	115.53	119.66
1	C	74	ASN	CB-CA-C	5.65	121.67	110.42
1	C	74	ASN	CA-C-O	5.46	128.32	120.51
1	C	339	ASP	CB-CA-C	5.42	118.66	109.55
1	C	481	GLY	CA-C-O	-5.41	109.45	120.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	82	ARG	CG-CD-NE	5.36	123.78	112.00
1	C	281	PHE	CB-CA-C	-5.34	101.91	110.14
1	D	321	GLU	CA-C-N	5.33	127.96	120.28
1	D	321	GLU	C-N-CA	5.33	127.96	120.28
1	C	257	GLN	N-CA-C	-5.30	102.21	110.10
1	B	74	ASN	N-CA-CB	-5.25	103.35	111.65
1	A	314	ASP	N-CA-C	5.24	119.42	113.19
1	A	80	ARG	CB-CG-CD	-5.23	99.28	111.30
1	B	445	ARG	CD-NE-CZ	5.22	131.71	124.40
1	A	-5	VAL	CA-C-N	-5.17	114.49	119.87
1	A	-5	VAL	C-N-CA	-5.17	114.49	119.87
1	D	445	ARG	CD-NE-CZ	5.12	131.56	124.40
1	A	426	ASP	N-CA-C	5.08	118.74	112.54
1	A	194	VAL	CB-CA-C	-5.05	101.98	110.71

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	3884	0	3729	6	0
1	B	3850	0	3689	5	0
1	C	3799	0	3633	9	0
1	D	3815	0	3655	5	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	268	0	0	1	0
3	B	296	0	0	1	0
3	C	233	0	0	3	0
3	D	279	0	0	0	0
All	All	16428	0	14706	25	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 1.

All (25) 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:405:ASP:OD2	3:C:601:HOH:O	2.15	0.65
1:A:272[A]:ARG:NH2	3:A:601:HOH:O	2.27	0.61
1:C:455:ASP:OD2	1:C:458:SER:OG	2.22	0.51
1:A:125:ILE:HD11	1:A:171:LEU:HG	1.93	0.51
1:C:225:GLU:OE1	3:C:602:HOH:O	2.20	0.50
1:A:3:ILE:O	1:A:52:PRO:HG2	2.12	0.49
1:D:281[B]:PHE:HZ	1:D:307:PHE:CE2	2.32	0.47
1:C:125:ILE:HD11	1:C:171:LEU:HG	1.96	0.47
1:D:281[B]:PHE:HZ	1:D:307:PHE:CD2	2.32	0.47
1:C:431:ASP:OD2	3:C:603:HOH:O	2.20	0.47
1:D:125:ILE:HD11	1:D:171:LEU:HG	1.96	0.47
1:B:428:GLU:HG3	3:B:860:HOH:O	2.17	0.45
1:B:124:VAL:HA	1:B:132:LEU:O	2.16	0.44
1:D:124:VAL:HA	1:D:132:LEU:O	2.17	0.44
1:D:233:ILE:HG23	1:D:281[A]:PHE:HB2	1.99	0.44
1:A:124:VAL:HA	1:A:132:LEU:O	2.18	0.43
1:B:125:ILE:HD11	1:B:171:LEU:HG	2.01	0.42
1:C:401:THR:O	1:C:403:LEU:CD2	2.67	0.42
1:A:-5:VAL:N	1:A:-4:PRO:CD	2.82	0.42
1:C:124:VAL:HA	1:C:132:LEU:O	2.20	0.42
1:B:233:ILE:HG23	1:B:281:PHE:HB2	2.02	0.42
1:C:306:ILE:HD13	1:C:306:ILE:HG21	1.78	0.41
1:C:135:LYS:HG3	1:C:137:ASP:OD1	2.20	0.41
1:A:272[B]:ARG:HA	1:A:272[B]:ARG:HD3	1.90	0.40
1:B:135:LYS:HG3	1:B:137:ASP:OD1	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	489/501 (98%)	483 (99%)	6 (1%)	0	100	100
1	B	485/501 (97%)	475 (98%)	9 (2%)	1 (0%)	43	36
1	C	478/501 (95%)	468 (98%)	8 (2%)	2 (0%)	30	22
1	D	480/501 (96%)	472 (98%)	8 (2%)	0	100	100
All	All	1932/2004 (96%)	1898 (98%)	31 (2%)	3 (0%)	43	36

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	314	ASP
1	C	314	ASP
1	C	74	ASN

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	411/419 (98%)	399 (97%)	12 (3%)	37	31
1	B	408/419 (97%)	398 (98%)	10 (2%)	42	37
1	C	401/419 (96%)	387 (96%)	14 (4%)	32	24
1	D	403/419 (96%)	391 (97%)	12 (3%)	36	30
All	All	1623/1676 (97%)	1575 (97%)	48 (3%)	37	30

All (48) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	-6	LEU
1	A	112	LEU
1	A	255	LEU
1	A	334	LEU
1	A	336	SER
1	A	338	ARG
1	A	363	ARG
1	A	364	GLN

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Mol	Chain	Res	Type
1	A	394	LEU
1	A	426	ASP
1	A	428	GLU
1	A	440	LEU
1	B	24[A]	GLU
1	B	24[B]	GLU
1	B	112	LEU
1	B	115	LYS
1	B	255	LEU
1	B	321	GLU
1	B	394	LEU
1	B	426	ASP
1	B	428	GLU
1	B	440	LEU
1	C	2	SER
1	C	34	LEU
1	C	74	ASN
1	C	112	LEU
1	C	255	LEU
1	C	328	MET
1	C	334	LEU
1	C	363	ARG
1	C	394	LEU
1	C	400	LYS
1	C	403	LEU
1	C	426	ASP
1	C	428	GLU
1	C	440	LEU
1	D	24	GLU
1	D	25	VAL
1	D	82[A]	ARG
1	D	82[B]	ARG
1	D	112	LEU
1	D	203	LYS
1	D	255	LEU
1	D	340	GLU
1	D	363	ARG
1	D	394	LEU
1	D	428	GLU
1	D	440	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (17) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	GLN
1	A	100	ASN
1	A	389	GLN
1	B	7	GLN
1	B	44	GLN
1	B	92	GLN
1	B	100	ASN
1	B	269	GLN
1	B	389	GLN
1	C	44	GLN
1	C	92	GLN
1	C	100	ASN
1	C	389	GLN
1	D	92	GLN
1	D	100	ASN
1	D	312	GLN
1	D	389	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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

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

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	488/501 (97%)	-0.18	5 (1%) 79 82	17, 32, 53, 93	3 (0%)
1	B	483/501 (96%)	-0.21	7 (1%) 73 76	17, 31, 51, 87	4 (0%)
1	C	480/501 (95%)	-0.00	13 (2%) 56 60	22, 34, 63, 95	0
1	D	480/501 (95%)	-0.23	4 (0%) 82 85	16, 31, 49, 70	2 (0%)
All	All	1931/2004 (96%)	-0.16	29 (1%) 72 75	16, 32, 54, 95	9 (0%)

All (29) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	388	PHE	3.6
1	C	386	ILE	3.3
1	B	306	ILE	3.3
1	C	379	TYR	2.7
1	D	2	SER	2.7
1	B	320	PRO	2.6
1	C	442	ALA	2.6
1	A	307	PHE	2.5
1	A	-6	LEU	2.5
1	C	381	PRO	2.5
1	B	307	PHE	2.5
1	D	481	GLY	2.4
1	C	3	ILE	2.4
1	C	2	SER	2.4
1	A	481	GLY	2.3
1	D	322	THR	2.3
1	B	322	THR	2.3
1	C	385	PRO	2.3
1	B	0[A]	HIS	2.2
1	A	338	ARG	2.2
1	C	382	ALA	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	308	TYR	2.1
1	A	306	ILE	2.1
1	C	441	ILE	2.1
1	C	403	LEU	2.1
1	D	307	PHE	2.0
1	C	401	THR	2.0
1	B	316	HIS	2.0
1	B	386	ILE	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(Å ²)	Q<0.9
2	FE	A	500	1/1	1.00	0.01	21,21,21,21	0
2	FE	B	500	1/1	1.00	0.01	21,21,21,21	0
2	FE	C	500	1/1	1.00	0.04	24,24,24,24	0
2	FE	D	500	1/1	1.00	0.02	21,21,21,21	0

6.5 Other polymers [i](#)

There are no such residues in this entry.