



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 11:09 PM UTC

PDB ID : 3OIC / pdb_00003oic
Title : Crystal Structure of Enoyl-ACP Reductases III (FabL) from *B. subtilis* (apo form)
Authors : Kim, K.-H.; Ha, B.H.; Kim, S.J.; Hong, S.K.; Hwang, K.Y.; Kim, E.E.
Deposited on : 2010-08-19
Resolution : 2.20 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

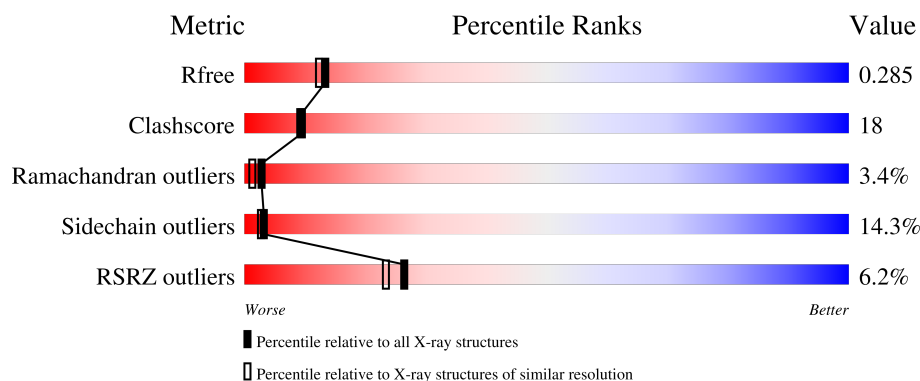
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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	6164 (2.20-2.20)
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)
RSRZ outliers	180081	6166 (2.20-2.20)

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	258	5% 65% 21% 7% • •
1	D	258	7% 66% 24% 5% • •

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3932 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 Enoyl-[acyl-carrier-protein] reductase [NADPH].

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	247	Total	C	N	O	S	0	0	0
			1889	1188	335	357	9			
1	D	252	Total	C	N	O	S	0	0	0
			1925	1210	340	366	9			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	251	LEU	-	expression tag	UNP P71079
A	252	GLU	-	expression tag	UNP P71079
A	253	HIS	-	expression tag	UNP P71079
A	254	HIS	-	expression tag	UNP P71079
A	255	HIS	-	expression tag	UNP P71079
A	256	HIS	-	expression tag	UNP P71079
A	257	HIS	-	expression tag	UNP P71079
A	258	HIS	-	expression tag	UNP P71079
D	251	LEU	-	expression tag	UNP P71079
D	252	GLU	-	expression tag	UNP P71079
D	253	HIS	-	expression tag	UNP P71079
D	254	HIS	-	expression tag	UNP P71079
D	255	HIS	-	expression tag	UNP P71079
D	256	HIS	-	expression tag	UNP P71079
D	257	HIS	-	expression tag	UNP P71079
D	258	HIS	-	expression tag	UNP P71079

- Molecule 2 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

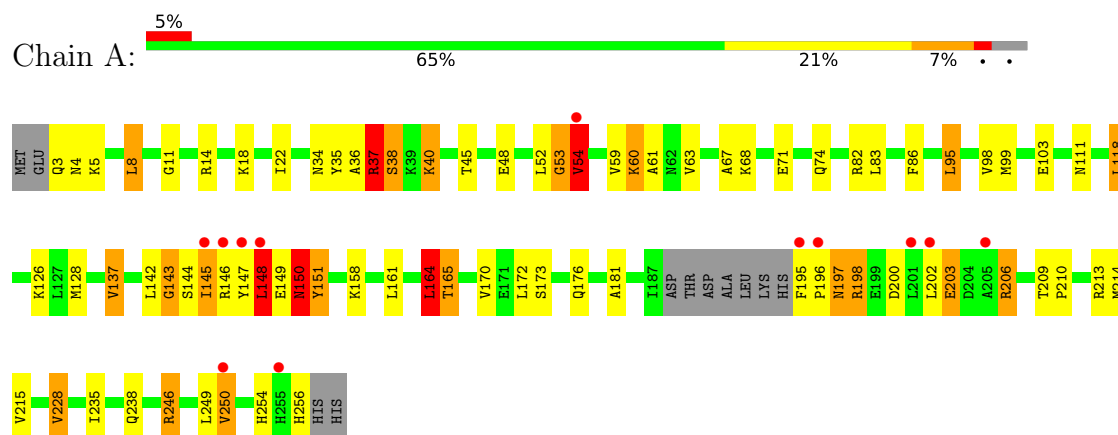
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	51	Total	O	0	0
			51	51		
3	D	62	Total	O	0	0
			62	62		

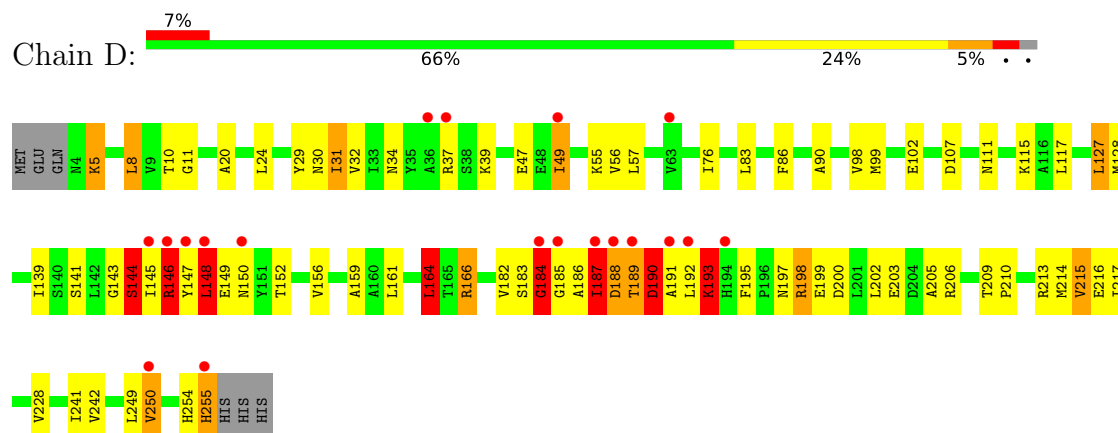
3 Residue-property plots [i](#)

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: Enoyl-[acyl-carrier-protein] reductase [NADPH]



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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	85.21Å 93.43Å 66.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.71 – 2.20 46.71 – 2.20	Depositor EDS
% Data completeness (in resolution range)	99.3 (46.71-2.20) 99.5 (46.71-2.20)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.20 (at 2.20Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.226 , 0.282 0.227 , 0.285	Depositor DCC
R_{free} test set	1388 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å ²)	34.1	Xtriage
Anisotropy	0.159	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 35.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	3932	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.18% 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: SO4

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	1.20	3/1913 (0.2%)	1.24	7/2581 (0.3%)
1	D	1.24	4/1950 (0.2%)	1.21	6/2633 (0.2%)
All	All	1.22	7/3863 (0.2%)	1.22	13/5214 (0.2%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	249	LEU	C-O	-5.87	1.16	1.24
1	D	159	ALA	CA-CB	-5.69	1.44	1.53
1	D	20	ALA	N-CA	5.53	1.53	1.46
1	A	63	VAL	CA-CB	5.53	1.63	1.54
1	D	128	MET	SD-CE	-5.52	1.65	1.79
1	A	228	VAL	C-O	-5.50	1.18	1.24
1	D	24	LEU	CA-C	5.28	1.59	1.52

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	215	VAL	CB-CA-C	-8.99	98.88	111.28
1	A	150	ASN	N-CA-C	-8.39	92.92	110.80
1	A	4	ASN	N-CA-C	8.21	121.51	110.35
1	A	53	GLY	N-CA-C	-7.40	96.91	112.45
1	D	188	ASP	N-CA-C	-7.29	104.64	113.97
1	D	184	GLY	N-CA-C	-7.12	96.30	113.18
1	D	187	ILE	CB-CA-C	-7.06	102.42	110.96
1	D	183	SER	N-CA-C	-5.95	106.05	113.55
1	A	164	LEU	CA-CB-CG	5.84	136.73	116.30
1	A	144	SER	N-CA-C	5.75	117.39	109.29
1	A	137	VAL	CB-CA-C	5.74	118.92	110.77
1	D	164	LEU	CA-CB-CG	5.31	134.89	116.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	151	TYR	CB-CA-C	5.13	120.63	110.42

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	1889	0	1930	61	0
1	D	1925	0	1967	83	0
2	D	5	0	0	0	0
3	A	51	0	0	7	0
3	D	62	0	0	6	0
All	All	3932	0	3897	138	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 18.

All (138) 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:D:189:THR:CG2	1:D:214:MET:HE3	1.78	1.14
1:D:189:THR:HG21	3:D:273:HOH:O	1.53	1.09
1:D:189:THR:HG22	1:D:214:MET:HE3	1.35	1.04
1:A:250:VAL:HG12	1:D:111:ASN:HD22	1.29	0.96
1:D:184:GLY:CA	1:D:242:VAL:H	1.80	0.94
1:D:184:GLY:HA2	1:D:242:VAL:H	1.29	0.94
1:A:250:VAL:HG12	1:D:111:ASN:ND2	1.85	0.92
1:A:147:TYR:HH	1:A:195:PHE:HZ	1.01	0.90
1:A:36:ALA:O	3:A:280:HOH:O	1.90	0.88
1:A:98:VAL:H	1:A:150:ASN:HA	1.40	0.85
1:A:60:LYS:HE3	1:A:68:LYS:HE2	1.61	0.81
1:D:49:ILE:HG13	1:D:56:VAL:HG21	1.62	0.81
1:A:37:ARG:O	1:A:38:SER:CB	2.31	0.79
1:D:209:THR:HG21	1:D:214:MET:SD	2.23	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:184:GLY:HA2	1:D:242:VAL:N	1.98	0.78
1:D:141:SER:HB2	1:D:186:ALA:CB	2.13	0.77
1:D:141:SER:HB2	1:D:186:ALA:HB3	1.65	0.77
1:D:189:THR:HG23	1:D:214:MET:HE3	1.66	0.75
1:D:184:GLY:C	1:D:186:ALA:H	1.94	0.75
1:D:189:THR:O	1:D:190:ASP:HB2	1.85	0.75
1:D:141:SER:CB	1:D:186:ALA:HB3	2.16	0.75
1:A:111:ASN:OD1	1:D:250:VAL:HG12	1.86	0.74
1:D:189:THR:HG22	1:D:214:MET:CE	2.14	0.73
3:A:298:HOH:O	1:D:185:GLY:HA3	1.88	0.73
1:D:205:ALA:O	1:D:209:THR:HG22	1.87	0.73
1:A:118:LEU:HD13	1:A:164:LEU:HG	1.72	0.71
1:A:40:LYS:HE3	1:A:40:LYS:HA	1.74	0.69
1:D:193:LYS:CB	1:D:193:LYS:NZ	2.56	0.69
1:A:256:HIS:CD2	3:A:267:HOH:O	2.46	0.69
1:A:53:GLY:O	1:A:54:VAL:HB	1.92	0.69
1:A:149:GLU:OE1	1:A:149:GLU:C	2.36	0.69
1:D:184:GLY:O	1:D:186:ALA:N	2.26	0.67
1:A:37:ARG:O	1:A:38:SER:HB3	1.95	0.66
1:A:142:LEU:O	1:A:143:GLY:O	2.14	0.66
1:D:190:ASP:O	1:D:193:LYS:HG2	1.96	0.66
1:A:36:ALA:O	1:A:37:ARG:O	2.14	0.65
1:D:193:LYS:HZ2	1:D:193:LYS:HB2	1.63	0.64
1:D:184:GLY:C	1:D:186:ALA:N	2.53	0.64
1:D:49:ILE:HG13	1:D:56:VAL:CG2	2.27	0.64
1:A:161:LEU:O	1:A:165:THR:CG2	2.46	0.64
1:D:98:VAL:HG22	1:D:150:ASN:HA	1.79	0.64
1:D:197:ASN:O	1:D:199:GLU:N	2.32	0.63
1:D:193:LYS:NZ	1:D:193:LYS:HB2	2.13	0.62
1:D:143:GLY:O	1:D:144:SER:C	2.42	0.62
1:A:246:ARG:HD2	1:D:107:ASP:HB2	1.81	0.61
1:A:165:THR:HB	3:A:273:HOH:O	2.01	0.61
1:D:141:SER:OG	1:D:186:ALA:HB3	2.00	0.60
1:A:11:GLY:H	1:A:34:ASN:HD22	1.49	0.59
1:D:189:THR:CG2	1:D:214:MET:CE	2.69	0.58
1:A:99:MET:HG3	3:D:308:HOH:O	2.04	0.58
1:A:161:LEU:O	1:A:165:THR:HG22	2.05	0.57
1:A:148:LEU:O	1:A:148:LEU:HD23	2.05	0.57
1:D:11:GLY:H	1:D:34:ASN:HD22	1.52	0.57
1:D:147:TYR:O	1:D:148:LEU:HB2	2.04	0.57
1:D:193:LYS:CB	1:D:193:LYS:HZ3	2.17	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:149:GLU:OE1	1:A:149:GLU:CA	2.53	0.56
1:D:8:LEU:HD13	1:D:86:PHE:CE1	2.40	0.56
1:D:166:ARG:HD3	3:D:313:HOH:O	2.04	0.56
1:A:197:ASN:O	1:A:200:ASP:N	2.39	0.55
1:A:149:GLU:O	1:A:149:GLU:CG	2.54	0.55
1:A:203:GLU:HG3	1:A:206:ARG:HH12	1.72	0.54
1:D:189:THR:CG2	3:D:273:HOH:O	2.30	0.54
1:D:32:VAL:HG21	1:D:76:ILE:HG12	1.90	0.54
1:D:146:ARG:HB3	1:D:195:PHE:HZ	1.73	0.54
1:D:199:GLU:O	1:D:203:GLU:HG2	2.07	0.54
1:D:29:TYR:O	1:D:31:ILE:HD13	2.08	0.53
1:D:254:HIS:O	1:D:255:HIS:HB2	2.08	0.53
1:A:8:LEU:HD13	1:A:86:PHE:CE1	2.44	0.53
1:A:256:HIS:HD2	3:A:267:HOH:O	1.87	0.53
1:D:146:ARG:HG2	1:D:147:TYR:N	2.24	0.52
1:A:161:LEU:O	1:A:165:THR:HG23	2.09	0.52
1:A:95:LEU:HG	1:A:151:TYR:HB3	1.90	0.52
1:A:149:GLU:O	1:A:149:GLU:HG3	2.08	0.52
1:D:99:MET:O	1:D:102:GLU:HG2	2.11	0.51
1:D:161:LEU:HA	1:D:164:LEU:HD13	1.91	0.51
1:A:147:TYR:O	1:A:149:GLU:N	2.44	0.51
1:A:149:GLU:OE1	1:A:149:GLU:O	2.29	0.51
1:A:37:ARG:O	1:A:38:SER:HB2	2.11	0.50
1:D:188:ASP:C	1:D:189:THR:O	2.50	0.50
1:A:149:GLU:O	1:A:150:ASN:HB2	2.12	0.50
1:D:30:ASN:ND2	1:D:55:LYS:H	2.09	0.50
1:D:98:VAL:CG2	1:D:150:ASN:HA	2.41	0.50
1:D:184:GLY:H	1:D:241:ILE:HA	1.76	0.50
1:A:209:THR:HG21	1:A:214:MET:HG2	1.93	0.50
1:D:10:THR:HA	1:D:34:ASN:HB3	1.94	0.49
1:A:60:LYS:CE	1:A:68:LYS:HE2	2.35	0.49
1:D:193:LYS:HZ3	1:D:193:LYS:HB3	1.77	0.49
1:D:146:ARG:HB3	1:D:195:PHE:CZ	2.47	0.49
1:A:98:VAL:H	1:A:150:ASN:CA	2.18	0.49
1:D:37:ARG:NH1	3:D:315:HOH:O	2.46	0.48
1:A:83:LEU:O	1:A:128:MET:HG2	2.13	0.48
1:A:210:PRO:HB2	1:A:213:ARG:HD3	1.94	0.48
1:D:197:ASN:O	1:D:198:ARG:C	2.57	0.48
1:A:197:ASN:O	1:A:198:ARG:C	2.57	0.48
1:D:98:VAL:HG12	1:D:152:THR:HG23	1.95	0.47
1:D:115:LYS:HE3	1:D:115:LYS:HB2	1.66	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:249:LEU:HD23	1:D:249:LEU:HA	1.76	0.47
1:D:146:ARG:CG	1:D:147:TYR:H	2.27	0.46
1:D:202:LEU:HD23	1:D:206:ARG:NH2	2.31	0.46
1:D:146:ARG:HG2	1:D:147:TYR:H	1.81	0.46
1:D:189:THR:O	1:D:190:ASP:CB	2.59	0.46
1:A:246:ARG:HD2	1:D:107:ASP:CB	2.45	0.46
1:D:139:ILE:HA	1:D:182:VAL:O	2.16	0.46
1:D:184:GLY:O	1:D:185:GLY:C	2.58	0.45
1:D:143:GLY:O	1:D:145:ILE:N	2.50	0.45
1:D:210:PRO:HB2	1:D:213:ARG:HD3	1.98	0.45
1:A:203:GLU:HG3	1:A:206:ARG:NH1	2.31	0.45
1:D:187:ILE:H	1:D:187:ILE:HG13	1.25	0.45
1:D:141:SER:HB2	1:D:186:ALA:HB2	1.97	0.45
1:A:52:LEU:C	1:A:53:GLY:O	2.55	0.44
1:D:187:ILE:HG21	1:D:217:ILE:N	2.32	0.44
1:A:8:LEU:HD13	1:A:86:PHE:CD1	2.53	0.44
1:A:18:LYS:HE2	3:A:265:HOH:O	2.17	0.44
1:A:161:LEU:HA	1:A:164:LEU:HD13	1.99	0.44
1:A:35:TYR:CZ	1:A:60:LYS:HB3	2.52	0.44
1:A:98:VAL:N	1:A:150:ASN:HA	2.21	0.43
1:A:196:PRO:C	1:A:198:ARG:H	2.25	0.43
1:D:146:ARG:CG	1:D:147:TYR:N	2.82	0.43
1:D:197:ASN:O	1:D:200:ASP:N	2.52	0.43
1:A:149:GLU:OE1	1:A:149:GLU:HA	2.19	0.43
1:D:187:ILE:CG2	1:D:216:GLU:HA	2.49	0.43
1:D:187:ILE:HG22	1:D:216:GLU:HA	2.00	0.43
1:A:67:ALA:O	1:A:71:GLU:HG3	2.19	0.42
1:A:98:VAL:HB	1:A:150:ASN:HA	2.01	0.42
1:A:103:GLU:HG3	1:D:195:PHE:CE1	2.54	0.42
1:D:83:LEU:HB2	1:D:127:LEU:HD13	2.01	0.42
1:D:190:ASP:C	1:D:192:LEU:N	2.77	0.42
1:A:36:ALA:O	1:A:37:ARG:C	2.62	0.42
1:A:165:THR:HG21	1:A:181:ALA:HB2	2.02	0.42
1:D:55:LYS:NZ	3:D:271:HOH:O	2.52	0.42
1:A:60:LYS:O	1:A:61:ALA:HB2	2.20	0.41
3:A:298:HOH:O	1:D:185:GLY:CA	2.59	0.41
1:A:235:ILE:HG23	1:A:238:GLN:HG3	2.02	0.41
1:D:148:LEU:HD12	1:D:148:LEU:HA	1.89	0.41
1:D:5:LYS:HD2	1:D:5:LYS:N	2.36	0.41
1:D:90:ALA:HB2	1:D:117:LEU:HD21	2.04	0.40
1:A:165:THR:HG21	1:A:181:ALA:CB	2.52	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:HIS:C	1:A:256:HIS:H	2.29	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	243/258 (94%)	219 (90%)	16 (7%)	8 (3%)	3	1
1	D	250/258 (97%)	232 (93%)	9 (4%)	9 (4%)	2	1
All	All	493/516 (96%)	451 (92%)	25 (5%)	17 (3%)	3	1

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	37	ARG
1	A	38	SER
1	A	143	GLY
1	A	148	LEU
1	D	144	SER
1	D	146	ARG
1	D	148	LEU
1	D	184	GLY
1	D	193	LYS
1	D	198	ARG
1	A	150	ASN
1	A	198	ARG
1	A	54	VAL
1	D	189	THR
1	D	190	ASP
1	D	191	ALA
1	A	145	ILE

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	201/211 (95%)	165 (82%)	36 (18%)	2	1
1	D	205/211 (97%)	183 (89%)	22 (11%)	6	6
All	All	406/422 (96%)	348 (86%)	58 (14%)	3	3

All (58) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	5	LYS
1	A	8	LEU
1	A	14	ARG
1	A	22	ILE
1	A	37	ARG
1	A	40	LYS
1	A	45	THR
1	A	48	GLU
1	A	54	VAL
1	A	59	VAL
1	A	60	LYS
1	A	74	GLN
1	A	82	ARG
1	A	95	LEU
1	A	118	LEU
1	A	126	LYS
1	A	137	VAL
1	A	145	ILE
1	A	146	ARG
1	A	148	LEU
1	A	158	LYS
1	A	164	LEU
1	A	165	THR
1	A	170	VAL
1	A	172	LEU
1	A	173	SER

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Mol	Chain	Res	Type
1	A	176	GLN
1	A	197	ASN
1	A	202	LEU
1	A	203	GLU
1	A	206	ARG
1	A	215	VAL
1	A	228	VAL
1	A	246	ARG
1	A	250	VAL
1	D	5	LYS
1	D	8	LEU
1	D	31	ILE
1	D	39	LYS
1	D	47	GLU
1	D	49	ILE
1	D	57	LEU
1	D	127	LEU
1	D	144	SER
1	D	146	ARG
1	D	148	LEU
1	D	149	GLU
1	D	156	VAL
1	D	164	LEU
1	D	166	ARG
1	D	187	ILE
1	D	190	ASP
1	D	193	LYS
1	D	215	VAL
1	D	228	VAL
1	D	250	VAL
1	D	255	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	3	GLN
1	A	30	ASN
1	A	34	ASN
1	A	88	ASN
1	A	256	HIS
1	D	27	ASN
1	D	30	ASN

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Mol	Chain	Res	Type
1	D	34	ASN
1	D	111	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	D	259	-	4,4,4	0.41	0	6,6,6	0.39	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 [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	247/258 (95%)	0.42	12 (4%) 35 31	21, 40, 64, 85	0
1	D	252/258 (97%)	0.28	19 (7%) 20 17	18, 33, 64, 74	0
All	All	499/516 (96%)	0.35	31 (6%) 26 23	18, 37, 64, 85	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	188	ASP	5.1
1	A	147	TYR	5.1
1	D	185	GLY	4.5
1	A	195	PHE	4.5
1	D	192	LEU	4.2
1	D	184	GLY	3.8
1	D	189	THR	3.7
1	D	191	ALA	3.5
1	D	187	ILE	3.4
1	A	145	ILE	3.2
1	D	49	ILE	3.2
1	D	145	ILE	3.0
1	A	250	VAL	3.0
1	D	194	HIS	3.0
1	A	255	HIS	2.9
1	D	147	TYR	2.8
1	D	37	ARG	2.7
1	A	54	VAL	2.7
1	A	196	PRO	2.6
1	D	36	ALA	2.5
1	A	202	LEU	2.5
1	D	148	LEU	2.4
1	D	255	HIS	2.3
1	A	148	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	146	ARG	2.2
1	D	150	ASN	2.2
1	A	205	ALA	2.2
1	D	146	ARG	2.2
1	D	63	VAL	2.1
1	D	250	VAL	2.0
1	A	201	LEU	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	SO4	D	259	5/5	0.78	0.12	62,66,68,70	0

6.5 Other polymers [i](#)

There are no such residues in this entry.