Full wwPDB X-ray Structure Validation Report (i)

Jun 16, 2024 – 02:39 PM EDT

PDB ID	:	4XDV
Title	:	Crystal Structure of the L74F/M78V/I80V/L114F mutant of LEH complexed with cyclohexanediol
Authors	:	Kong, X.D.; Sun, Z.; Lonsdale, R.; Xu, J.H.; Reetz, M.T.; Zhou, J.
Deposited on	:	2014-12-20
Resolution	:	2.25 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 (i) 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 (i)) were used in the production of this report:

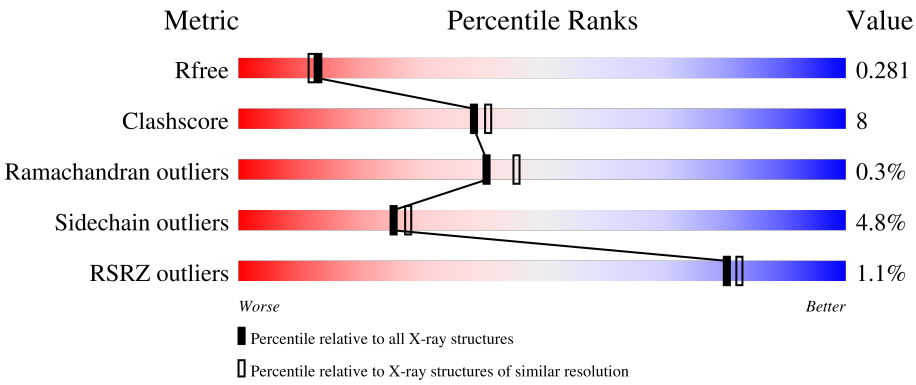
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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}	130704	1377 (2.26-2.26)
Clashscore	141614	1487 (2.26-2.26)
Ramachandran outliers	138981	1449 (2.26-2.26)
Sidechain outliers	138945	1450 (2.26-2.26)
RSRZ outliers	127900	1356 (2.26-2.26)

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 <=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	155	.% 81% 10% • 8%
1	B	155	79% 13% • 6%
1	C	155	.% 77% 15% 7%
1	D	155	81% 12% • 6%
1	E	155	3% 57% 21% • 19%

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Mol	Chain	Length	Quality of chain
1	F	155	3% 74% 14% • 10%
1	G	155	% 76% 15% • 8%
1	H	155	% 71% 15% • • 9%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 9569 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 Limonene-1,2-epoxide hydrolase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	C	144	1132	724	182	223	3	0	0	0
			Total	C	N	O	S			
1	E	126	989	629	158	199	3	0	0	0
			Total	C	N	O	S			
1	G	143	1121	718	178	222	3	0	0	0
			Total	C	N	O	S			
1	D	145	1136	726	183	224	3	0	0	0
			Total	C	N	O	S			
1	F	140	1098	701	175	219	3	0	0	0
			Total	C	N	O	S			
1	H	141	1106	707	176	220	3	0	0	0
			Total	C	N	O	S			
1	A	143	1132	727	179	223	3	0	1	0
			Total	C	N	O	S			
1	B	145	1147	735	184	225	3	0	1	0

There are 112 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	-5	MET	-	initiating methionine	UNP Q9ZAG3
C	-4	HIS	-	expression tag	UNP Q9ZAG3
C	-3	HIS	-	expression tag	UNP Q9ZAG3
C	-2	HIS	-	expression tag	UNP Q9ZAG3
C	-1	HIS	-	expression tag	UNP Q9ZAG3
C	0	HIS	-	expression tag	UNP Q9ZAG3
C	1	HIS	-	expression tag	UNP Q9ZAG3
C	2	THR	-	expression tag	UNP Q9ZAG3
C	3	SER	-	expression tag	UNP Q9ZAG3
C	4	LEU	-	expression tag	UNP Q9ZAG3
C	74	PHE	LEU	engineered mutation	UNP Q9ZAG3
C	78	VAL	MET	engineered mutation	UNP Q9ZAG3
C	80	VAL	ILE	engineered mutation	UNP Q9ZAG3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	114	PHE	LEU	engineered mutation	UNP Q9ZAG3
E	-5	MET	-	initiating methionine	UNP Q9ZAG3
E	-4	HIS	-	expression tag	UNP Q9ZAG3
E	-3	HIS	-	expression tag	UNP Q9ZAG3
E	-2	HIS	-	expression tag	UNP Q9ZAG3
E	-1	HIS	-	expression tag	UNP Q9ZAG3
E	0	HIS	-	expression tag	UNP Q9ZAG3
E	1	HIS	-	expression tag	UNP Q9ZAG3
E	2	THR	-	expression tag	UNP Q9ZAG3
E	3	SER	-	expression tag	UNP Q9ZAG3
E	4	LEU	-	expression tag	UNP Q9ZAG3
E	74	PHE	LEU	engineered mutation	UNP Q9ZAG3
E	78	VAL	MET	engineered mutation	UNP Q9ZAG3
E	80	VAL	ILE	engineered mutation	UNP Q9ZAG3
E	114	PHE	LEU	engineered mutation	UNP Q9ZAG3
G	-5	MET	-	initiating methionine	UNP Q9ZAG3
G	-4	HIS	-	expression tag	UNP Q9ZAG3
G	-3	HIS	-	expression tag	UNP Q9ZAG3
G	-2	HIS	-	expression tag	UNP Q9ZAG3
G	-1	HIS	-	expression tag	UNP Q9ZAG3
G	0	HIS	-	expression tag	UNP Q9ZAG3
G	1	HIS	-	expression tag	UNP Q9ZAG3
G	2	THR	-	expression tag	UNP Q9ZAG3
G	3	SER	-	expression tag	UNP Q9ZAG3
G	4	LEU	-	expression tag	UNP Q9ZAG3
G	74	PHE	LEU	engineered mutation	UNP Q9ZAG3
G	78	VAL	MET	engineered mutation	UNP Q9ZAG3
G	80	VAL	ILE	engineered mutation	UNP Q9ZAG3
G	114	PHE	LEU	engineered mutation	UNP Q9ZAG3
D	-5	MET	-	initiating methionine	UNP Q9ZAG3
D	-4	HIS	-	expression tag	UNP Q9ZAG3
D	-3	HIS	-	expression tag	UNP Q9ZAG3
D	-2	HIS	-	expression tag	UNP Q9ZAG3
D	-1	HIS	-	expression tag	UNP Q9ZAG3
D	0	HIS	-	expression tag	UNP Q9ZAG3
D	1	HIS	-	expression tag	UNP Q9ZAG3
D	2	THR	-	expression tag	UNP Q9ZAG3
D	3	SER	-	expression tag	UNP Q9ZAG3
D	4	LEU	-	expression tag	UNP Q9ZAG3
D	74	PHE	LEU	engineered mutation	UNP Q9ZAG3
D	78	VAL	MET	engineered mutation	UNP Q9ZAG3
D	80	VAL	ILE	engineered mutation	UNP Q9ZAG3

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Chain	Residue	Modelled	Actual	Comment	Reference
D	114	PHE	LEU	engineered mutation	UNP Q9ZAG3
F	-5	MET	-	initiating methionine	UNP Q9ZAG3
F	-4	HIS	-	expression tag	UNP Q9ZAG3
F	-3	HIS	-	expression tag	UNP Q9ZAG3
F	-2	HIS	-	expression tag	UNP Q9ZAG3
F	-1	HIS	-	expression tag	UNP Q9ZAG3
F	0	HIS	-	expression tag	UNP Q9ZAG3
F	1	HIS	-	expression tag	UNP Q9ZAG3
F	2	THR	-	expression tag	UNP Q9ZAG3
F	3	SER	-	expression tag	UNP Q9ZAG3
F	4	LEU	-	expression tag	UNP Q9ZAG3
F	74	PHE	LEU	engineered mutation	UNP Q9ZAG3
F	78	VAL	MET	engineered mutation	UNP Q9ZAG3
F	80	VAL	ILE	engineered mutation	UNP Q9ZAG3
F	114	PHE	LEU	engineered mutation	UNP Q9ZAG3
H	-5	MET	-	initiating methionine	UNP Q9ZAG3
H	-4	HIS	-	expression tag	UNP Q9ZAG3
H	-3	HIS	-	expression tag	UNP Q9ZAG3
H	-2	HIS	-	expression tag	UNP Q9ZAG3
H	-1	HIS	-	expression tag	UNP Q9ZAG3
H	0	HIS	-	expression tag	UNP Q9ZAG3
H	1	HIS	-	expression tag	UNP Q9ZAG3
H	2	THR	-	expression tag	UNP Q9ZAG3
H	3	SER	-	expression tag	UNP Q9ZAG3
H	4	LEU	-	expression tag	UNP Q9ZAG3
H	74	PHE	LEU	engineered mutation	UNP Q9ZAG3
H	78	VAL	MET	engineered mutation	UNP Q9ZAG3
H	80	VAL	ILE	engineered mutation	UNP Q9ZAG3
H	114	PHE	LEU	engineered mutation	UNP Q9ZAG3
A	-5	MET	-	initiating methionine	UNP Q9ZAG3
A	-4	HIS	-	expression tag	UNP Q9ZAG3
A	-3	HIS	-	expression tag	UNP Q9ZAG3
A	-2	HIS	-	expression tag	UNP Q9ZAG3
A	-1	HIS	-	expression tag	UNP Q9ZAG3
A	0	HIS	-	expression tag	UNP Q9ZAG3
A	1	HIS	-	expression tag	UNP Q9ZAG3
A	2	THR	-	expression tag	UNP Q9ZAG3
A	3	SER	-	expression tag	UNP Q9ZAG3
A	4	LEU	-	expression tag	UNP Q9ZAG3
A	74	PHE	LEU	engineered mutation	UNP Q9ZAG3
A	78	VAL	MET	engineered mutation	UNP Q9ZAG3
A	80	VAL	ILE	engineered mutation	UNP Q9ZAG3

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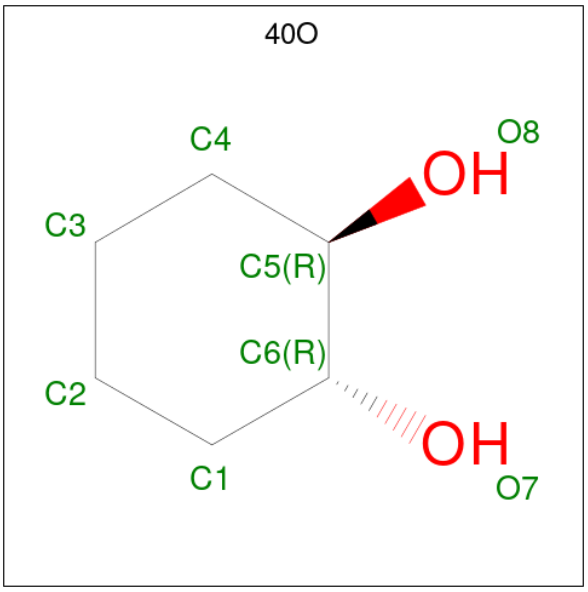
Chain	Residue	Modelled	Actual	Comment	Reference
A	114	PHE	LEU	engineered mutation	UNP Q9ZAG3
B	-5	MET	-	initiating methionine	UNP Q9ZAG3
B	-4	HIS	-	expression tag	UNP Q9ZAG3
B	-3	HIS	-	expression tag	UNP Q9ZAG3
B	-2	HIS	-	expression tag	UNP Q9ZAG3
B	-1	HIS	-	expression tag	UNP Q9ZAG3
B	0	HIS	-	expression tag	UNP Q9ZAG3
B	1	HIS	-	expression tag	UNP Q9ZAG3
B	2	THR	-	expression tag	UNP Q9ZAG3
B	3	SER	-	expression tag	UNP Q9ZAG3
B	4	LEU	-	expression tag	UNP Q9ZAG3
B	74	PHE	LEU	engineered mutation	UNP Q9ZAG3
B	78	VAL	MET	engineered mutation	UNP Q9ZAG3
B	80	VAL	ILE	engineered mutation	UNP Q9ZAG3
B	114	PHE	LEU	engineered mutation	UNP Q9ZAG3

- Molecule 2 is (1R,2R)-cyclohexane-1,2-diol (three-letter code: 40O) (formula: $C_6H_{12}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	C	1	8	6	2	0	0
2	D	1	8	6	2	1	0
2	A	1	8	6	2	0	0
2	B	1	8	6	2	1	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	C	96	Total O 96 96	0	0
3	E	58	Total O 58 58	0	0
3	G	63	Total O 63 63	0	0
3	D	93	Total O 93 93	0	0
3	F	52	Total O 52 52	0	0
3	H	67	Total O 67 67	0	0
3	A	120	Total O 120 120	0	0
3	B	127	Total O 127 127	0	0

3 Residue-property plots (i)

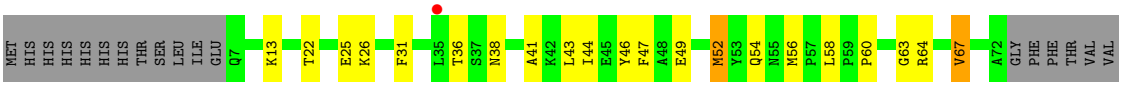
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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.

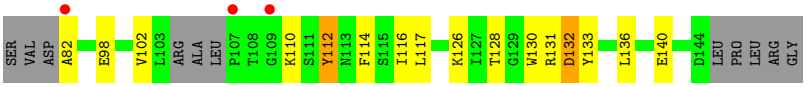
• Molecule 1: Limonene-1,2-epoxide hydrolase

Chain C:

• Molecule 1: Limonene-1,2-epoxide hydrolase

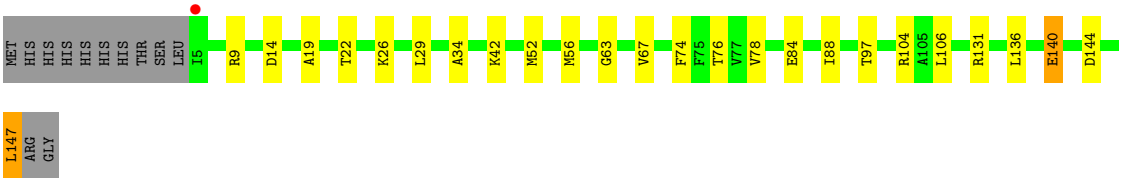
Chain E:





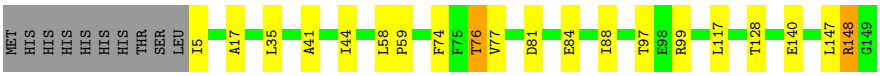
• Molecule 1: Limonene-1,2-epoxide hydrolase

Chain G:



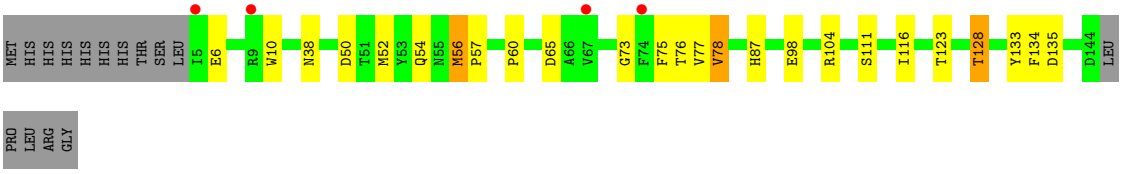
• Molecule 1: Limonene-1,2-epoxide hydrolase

Chain D:



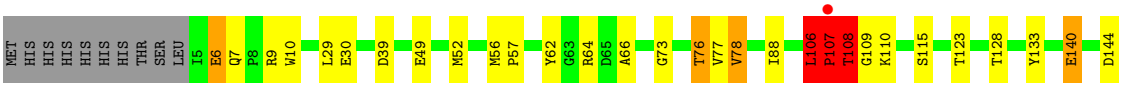
• Molecule 1: Limonene-1,2-epoxide hydrolase

Chain F:



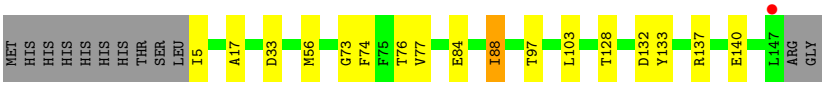
• Molecule 1: Limonene-1,2-epoxide hydrolase

Chain H:



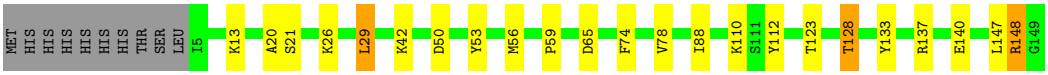
• Molecule 1: Limonene-1,2-epoxide hydrolase

Chain A:



• Molecule 1: Limonene-1,2-epoxide hydrolase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	104.70Å 61.16Å 119.55Å 90.00° 90.15° 90.00°	Depositor
Resolution (Å)	33.52 – 2.25 33.52 – 2.24	Depositor EDS
% Data completeness (in resolution range)	96.6 (33.52-2.25) 96.6 (33.52-2.24)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.27 (at 2.24Å)	Xtriage
Refinement program	PHENIX	Depositor
R, R_{free}	0.221 , 0.281 0.222 , 0.281	Depositor DCC
R_{free} test set	3505 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å^2)	39.4	Xtriage
Anisotropy	0.359	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 21.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.47$, $< L^2 > = 0.30$	Xtriage
Estimated twinning fraction	0.438 for h,-k,-l	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	9569	wwPDB-VP
Average B, all atoms (Å^2)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 61.82 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 1.2374e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: 4OO

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	0.47	0/1159	0.58	0/1577
1	B	0.46	0/1174	0.57	0/1596
1	C	0.47	0/1158	0.56	0/1575
1	D	0.44	0/1162	0.54	0/1580
1	E	0.33	0/1010	0.49	0/1369
1	F	0.34	0/1123	0.48	0/1527
1	G	0.35	0/1147	0.51	0/1561
1	H	0.36	0/1131	0.57	1/1538 (0.1%)
All	All	0.41	0/9064	0.54	1/12323 (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	H	0	2

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	H	108	THR	N-CA-C	-5.84	95.24	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	H	106	LEU	Peptide

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Mol	Chain	Res	Type	Group
1	H	107	PRO	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	1132	0	1085	12	0
1	B	1147	0	1101	19	0
1	C	1132	0	1090	17	0
1	D	1136	0	1093	16	0
1	E	989	0	937	26	0
1	F	1098	0	1048	15	0
1	G	1121	0	1077	14	1
1	H	1106	0	1059	18	0
2	A	8	0	12	3	0
2	B	8	0	12	4	0
2	C	8	0	12	3	0
2	D	8	0	12	1	0
3	A	120	0	0	5	0
3	B	127	0	0	4	0
3	C	96	0	0	5	0
3	D	93	0	0	2	0
3	E	58	0	0	5	0
3	F	52	0	0	1	0
3	G	63	0	0	1	0
3	H	67	0	0	1	0
All	All	9569	0	8538	131	1

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 8.

All (131) 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)
2:A:201:40O:O8	3:A:420:HOH:O	1.82	0.87
1:D:74:PHE:HB3	2:D:201:40O:H9	1.66	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:44:ILE:O	1:E:64:ARG:NH1	2.17	0.77
1:C:118:GLY:O	3:C:369:HOH:O	2.03	0.76
1:H:6:GLU:HB3	1:H:7:GLN:HA	1.67	0.76
1:A:132:ASP:OD2	3:A:420:HOH:O	2.05	0.74
1:A:74[B]:PHE:CZ	2:A:201:4OO:H1	2.23	0.74
1:E:49:GLU:OE2	1:E:64:ARG:HD3	1.89	0.72
1:E:54:GLN:HB2	1:E:60:PRO:HB3	1.71	0.71
1:C:53:TYR:OH	2:C:201:4OO:H2	1.90	0.70
1:E:130:TRP:NE1	1:E:132:ASP:OD1	2.25	0.70
1:A:140:GLU:OE2	3:A:414:HOH:O	2.08	0.70
1:C:148:ARG:NH2	3:C:380:HOH:O	2.24	0.68
1:E:43:LEU:O	3:E:231:HOH:O	2.13	0.67
1:F:65:ASP:OD2	3:F:201:HOH:O	2.12	0.67
1:H:6:GLU:HB3	1:H:7:GLN:CA	2.24	0.67
1:B:59:PRO:HG2	1:B:148:ARG:HB3	1.74	0.67
1:C:113:ASN:O	3:C:373:HOH:O	2.12	0.66
1:D:76:THR:HG22	1:D:77:VAL:HG13	1.79	0.65
1:B:20:ALA:O	3:B:379:HOH:O	2.15	0.64
1:H:108:THR:O	1:H:108:THR:OG1	2.14	0.63
1:A:88:ILE:HD12	1:A:97:THR:HG22	1.81	0.62
1:D:59:PRO:HG2	1:D:148:ARG:HB3	1.82	0.61
1:H:106:LEU:HD12	1:H:107:PRO:HD3	1.83	0.61
1:G:104:ARG:HH11	1:G:104:ARG:HG2	1.65	0.60
1:E:26:LYS:O	3:E:255:HOH:O	2.16	0.60
1:G:140:GLU:HG3	1:G:147:LEU:HD11	1.83	0.60
1:E:126:LYS:NZ	3:E:235:HOH:O	2.14	0.59
1:B:74[B]:PHE:CD2	2:B:201:4OO:H7	2.38	0.58
1:E:47:PHE:O	1:E:64:ARG:NH1	2.36	0.58
1:A:5:ILE:HG22	1:A:84:GLU:HG3	1.86	0.58
1:G:88:ILE:HD12	1:G:97:THR:HG22	1.84	0.58
1:H:108:THR:O	1:H:110:LYS:N	2.28	0.58
1:E:38:ASN:O	3:E:258:HOH:O	2.17	0.57
1:C:97:THR:N	3:C:369:HOH:O	2.37	0.57
1:F:123:THR:OG1	1:F:128:THR:HG21	2.05	0.56
1:F:104:ARG:HG2	1:F:104:ARG:HH11	1.70	0.56
1:F:77:VAL:HG23	1:F:78:VAL:HG12	1.87	0.56
1:B:42:LYS:NZ	3:B:372:HOH:O	2.19	0.55
1:D:88:ILE:HD12	1:D:97:THR:HG22	1.89	0.55
1:E:31:PHE:HE1	1:E:43:LEU:HD22	1.72	0.54
1:D:81:ASP:OD2	3:D:301:HOH:O	2.18	0.54
1:C:59:PRO:HG2	1:C:148:ARG:HB2	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:75:PHE:CE1	2:C:201:40O:H8	2.43	0.54
1:H:9:ARG:HB3	1:H:9:ARG:NH1	2.23	0.53
1:B:53:TYR:OH	2:B:201:40O:H6	2.08	0.53
1:G:131:ARG:NH1	1:H:10:TRP:HB2	2.24	0.53
1:E:56:MET:HG3	1:E:133:TYR:CE2	2.45	0.52
1:E:22:THR:OG1	1:E:25:GLU:HG3	2.09	0.52
1:C:140:GLU:HB3	3:C:391:HOH:O	2.10	0.52
1:D:147:LEU:O	1:D:148:ARG:HB2	2.11	0.51
1:F:116:ILE:HG12	1:F:134:PHE:HB3	1.93	0.50
1:E:54:GLN:NE2	1:E:58:LEU:O	2.45	0.49
1:C:22:THR:O	1:C:26:LYS:HG2	2.12	0.49
1:B:123:THR:OG1	1:B:128:THR:HG21	2.11	0.49
1:E:63:GLY:O	1:E:67:VAL:HG13	2.11	0.49
1:E:116:ILE:HG21	1:E:132:ASP:OD2	2.13	0.49
1:A:73:GLY:O	1:A:76:THR:HB	2.13	0.49
1:D:140:GLU:HA	1:D:147:LEU:HD21	1.94	0.49
1:G:14:ASP:HB3	1:B:21:SER:HB2	1.95	0.48
1:F:54:GLN:HB2	1:F:60:PRO:HB3	1.94	0.48
1:E:52:MET:HE1	1:E:131:ARG:HH11	1.78	0.48
1:B:110:LYS:HD3	1:B:112:TYR:CZ	2.49	0.48
1:G:104:ARG:HG2	1:G:104:ARG:NH1	2.27	0.48
1:B:140:GLU:HB3	1:B:147:LEU:HD22	1.96	0.48
1:G:19:ALA:O	3:G:239:HOH:O	2.20	0.48
1:G:22:THR:O	1:G:26:LYS:HG2	2.14	0.48
1:G:56:MET:O	1:G:136:LEU:HD11	2.13	0.48
1:A:33:ASP:OD1	3:A:390:HOH:O	2.20	0.47
1:H:73:GLY:O	1:H:76:THR:HG22	2.14	0.47
3:A:414:HOH:O	1:B:137:ARG:NH1	2.43	0.47
1:H:29:LEU:HG	1:H:88:ILE:HD13	1.97	0.47
1:D:17:ALA:HA	1:D:88:ILE:HG22	1.97	0.47
1:A:17:ALA:HA	1:A:88:ILE:HG22	1.97	0.47
1:E:112:TYR:HB2	1:E:114:PHE:CE1	2.50	0.46
1:F:56:MET:HA	1:F:57:PRO:HA	1.76	0.46
1:B:74[B]:PHE:CE1	1:B:78:VAL:HG11	2.51	0.46
1:C:117:LEU:HD23	1:D:117:LEU:HG	1.96	0.46
1:D:58:LEU:HD22	1:D:147:LEU:HB3	1.98	0.46
1:A:76:THR:HG22	1:A:77:VAL:HG13	1.96	0.46
1:G:14:ASP:OD2	1:B:26:LYS:NZ	2.45	0.46
1:E:41:ALA:O	1:E:44:ILE:HG22	2.16	0.45
1:H:140:GLU:H	1:H:140:GLU:HG2	1.46	0.45
1:D:5:ILE:N	3:D:378:HOH:O	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:73:GLY:O	1:F:76:THR:HB	2.17	0.45
1:C:6:GLU:OE2	1:C:86:PHE:CZ	2.70	0.45
1:F:104:ARG:HG2	1:F:104:ARG:NH1	2.31	0.45
1:C:117:LEU:CD2	1:D:117:LEU:HG	2.46	0.44
1:H:106:LEU:HD12	1:H:106:LEU:HA	1.79	0.44
1:E:136:LEU:HD12	1:F:135:ASP:OD2	2.17	0.44
1:B:29:LEU:HD12	1:B:29:LEU:HA	1.90	0.44
1:D:41:ALA:O	1:D:44:ILE:HG22	2.16	0.44
2:B:201:40O:H3	3:B:427:HOH:O	2.18	0.44
1:H:62:TYR:O	1:H:66:ALA:HB3	2.18	0.44
1:A:74[B]:PHE:CE2	2:A:201:40O:H9	2.53	0.44
1:E:131:ARG:NH1	1:F:10:TRP:HB2	2.33	0.43
1:C:80:VAL:HG21	2:C:201:40O:H7	2.00	0.43
1:H:56:MET:HG3	1:H:133:TYR:CZ	2.53	0.43
1:A:137:ARG:NH1	1:B:147:LEU:HD23	2.33	0.43
1:E:117:LEU:HD12	1:E:117:LEU:HA	1.79	0.43
1:B:74[B]:PHE:CE2	2:B:201:40O:H7	2.53	0.43
1:D:148:ARG:N	1:D:148:ARG:HD2	2.34	0.43
1:C:23:PRO:O	1:C:27:ILE:HG13	2.19	0.43
1:A:56:MET:HG3	1:A:133:TYR:CZ	2.54	0.43
1:H:78:VAL:N	1:H:106:LEU:HD22	2.33	0.43
1:G:63:GLY:O	1:G:67:VAL:HG23	2.19	0.42
1:G:74:PHE:O	1:G:78:VAL:HG12	2.19	0.42
1:F:56:MET:HG2	1:F:133:TYR:CE2	2.54	0.42
1:H:56:MET:HA	1:H:57:PRO:HA	1.81	0.42
1:B:29:LEU:HD13	1:B:88:ILE:HD13	2.02	0.42
1:E:114:PHE:O	1:E:116:ILE:HG13	2.20	0.42
1:C:73:GLY:O	1:C:76:THR:HB	2.20	0.42
1:F:87:HIS:HB2	1:F:98:GLU:HB3	2.02	0.42
1:C:74:PHE:O	1:C:78:VAL:HG12	2.20	0.41
1:G:76:THR:O	1:G:106:LEU:HD12	2.20	0.41
1:H:49:GLU:HA	1:H:64:ARG:HB2	2.02	0.41
1:D:35:LEU:HA	1:D:35:LEU:HD23	1.83	0.41
1:E:98:GLU:HG3	1:E:117:LEU:HD13	2.03	0.41
1:C:106:LEU:HB2	1:C:107:PRO:HD3	2.02	0.41
1:H:39:ASP:HB2	3:H:265:HOH:O	2.20	0.41
1:E:46:TYR:HB2	3:E:231:HOH:O	2.21	0.41
1:F:38:ASN:HA	1:F:75:PHE:CE1	2.55	0.41
1:B:56:MET:HG3	1:B:133:TYR:CZ	2.56	0.41
1:E:31:PHE:CE1	1:E:43:LEU:HD22	2.54	0.41
1:G:34:ALA:HB1	1:G:42:LYS:HE3	2.02	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:116:ILE:HG12	1:F:134:PHE:CB	2.51	0.40
1:H:123:THR:OG1	1:H:128:THR:HG21	2.21	0.40
1:E:82:ALA:HB3	1:E:102:VAL:HB	2.03	0.40
1:D:84:GLU:O	1:D:99:ARG:HB2	2.22	0.40
1:B:65:ASP:OD2	3:B:301:HOH:O	2.22	0.40
1:B:147:LEU:O	1:B:148:ARG:HB2	2.20	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:9:ARG:NH1	1:G:144:ASP:OD2[2_857]	2.00	0.20

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	142/155 (92%)	140 (99%)	2 (1%)	0	100	100
1	B	144/155 (93%)	141 (98%)	3 (2%)	0	100	100
1	C	142/155 (92%)	137 (96%)	5 (4%)	0	100	100
1	D	143/155 (92%)	141 (99%)	2 (1%)	0	100	100
1	E	120/155 (77%)	111 (92%)	9 (8%)	0	100	100
1	F	138/155 (89%)	134 (97%)	4 (3%)	0	100	100
1	G	141/155 (91%)	139 (99%)	2 (1%)	0	100	100
1	H	139/155 (90%)	129 (93%)	7 (5%)	3 (2%)	6	3
All	All	1109/1240 (89%)	1072 (97%)	34 (3%)	3 (0%)	41	46

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	H	107	PRO
1	H	144	ASP
1	H	109	GLY

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	120/130 (92%)	117 (98%)	3 (2%)	47	56
1	B	121/130 (93%)	116 (96%)	5 (4%)	30	36
1	C	120/130 (92%)	118 (98%)	2 (2%)	60	71
1	D	120/130 (92%)	117 (98%)	3 (2%)	47	56
1	E	104/130 (80%)	95 (91%)	9 (9%)	10	8
1	F	116/130 (89%)	109 (94%)	7 (6%)	19	18
1	G	119/130 (92%)	114 (96%)	5 (4%)	30	34
1	H	117/130 (90%)	106 (91%)	11 (9%)	8	6
All	All	937/1040 (90%)	892 (95%)	45 (5%)	25	28

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	C	5	ILE
1	C	128	THR
1	E	13	LYS
1	E	36	THR
1	E	52	MET
1	E	67	VAL
1	E	110	LYS
1	E	112	TYR
1	E	128	THR
1	E	132	ASP
1	E	140	GLU
1	G	29	LEU
1	G	52	MET

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Mol	Chain	Res	Type
1	G	84	GLU
1	G	140	GLU
1	G	147	LEU
1	D	76	THR
1	D	128	THR
1	D	148	ARG
1	F	6	GLU
1	F	50	ASP
1	F	52	MET
1	F	56	MET
1	F	78	VAL
1	F	111	SER
1	F	128	THR
1	H	6	GLU
1	H	30	GLU
1	H	52	MET
1	H	76	THR
1	H	77	VAL
1	H	78	VAL
1	H	106	LEU
1	H	108	THR
1	H	115	SER
1	H	140	GLU
1	H	145	LEU
1	A	88	ILE
1	A	103	LEU
1	A	128	THR
1	B	13	LYS
1	B	29	LEU
1	B	50	ASP
1	B	128	THR
1	B	148	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

4 ligands 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	$\#\|Z\| > 2$
2	40O	D	201	-	8,8,8	0.52	0	10,10,10	1.57	2 (20%)
2	40O	B	201	-	8,8,8	0.40	0	10,10,10	1.85	3 (30%)
2	40O	A	201	-	8,8,8	0.41	0	10,10,10	2.28	2 (20%)
2	40O	C	201	-	8,8,8	0.50	0	10,10,10	1.52	1 (10%)

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	40O	D	201	-	-	-	0/1/1/1
2	40O	B	201	-	-	-	0/1/1/1
2	40O	A	201	-	-	-	0/1/1/1
2	40O	C	201	-	-	-	0/1/1/1

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	201	40O	C4-C5-C6	5.84	115.79	111.12
2	B	201	40O	C2-C1-C6	3.75	118.32	111.47
2	A	201	40O	C3-C4-C5	3.63	118.11	111.47
2	B	201	40O	C3-C2-C1	3.61	118.77	111.42
2	D	201	40O	C1-C6-C5	-3.05	108.69	111.12
2	C	201	40O	C1-C6-C5	2.93	113.46	111.12
2	D	201	40O	C2-C3-C4	2.72	116.97	111.42
2	B	201	40O	C4-C5-C6	-2.09	109.45	111.12

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	201	40O	1	0
2	B	201	40O	4	0
2	A	201	40O	3	0
2	C	201	40O	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	143/155 (92%)	-0.44	1 (0%)	87	88	13, 19, 35, 50	0
1	B	145/155 (93%)	-0.46	0	100	100	12, 20, 35, 56	0
1	C	144/155 (92%)	-0.38	1 (0%)	87	88	17, 25, 42, 51	0
1	D	145/155 (93%)	-0.44	0	100	100	16, 24, 39, 56	0
1	E	126/155 (81%)	-0.01	4 (3%)	47	50	27, 48, 79, 85	0
1	F	140/155 (90%)	-0.17	4 (2%)	51	55	28, 47, 61, 67	0
1	G	143/155 (92%)	-0.26	1 (0%)	87	88	20, 42, 60, 84	0
1	H	141/155 (90%)	-0.20	1 (0%)	87	88	21, 42, 63, 77	0
All	All	1127/1240 (90%)	-0.30	12 (1%)	80	82	12, 31, 61, 85	0

All (12) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	H	107	PRO	5.0
1	E	107	PRO	4.4
1	G	5	ILE	3.1
1	E	35	LEU	2.9
1	C	6	GLU	2.7
1	A	147	LEU	2.6
1	F	74	PHE	2.5
1	E	82	ALA	2.5
1	F	9	ARG	2.4
1	F	5	ILE	2.1
1	F	67	VAL	2.1
1	E	109	GLY	2.1

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 monosaccharides 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	4OO	A	201	8/8	0.85	0.32	26,33,37,37	0
2	4OO	D	201	8/8	0.88	0.21	27,29,32,32	4
2	4OO	B	201	8/8	0.90	0.39	21,23,30,34	5
2	4OO	C	201	8/8	0.91	0.22	26,27,31,33	6

6.5 Other polymers (i)

There are no such residues in this entry.