



Full wwPDB X-ray Structure Validation Report ⓘ

May 23, 2020 – 07:09 am BST

PDB ID : 5XV3
Title : Crystal structure of ATG101-ATG13HORMA
Authors : Kim, B.-W.; Song, H.K.
Deposited on : 2017-06-26
Resolution : 2.57 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 following versions of software and data (see [references ⓘ](#)) 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 : 2.11
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) : 2.11

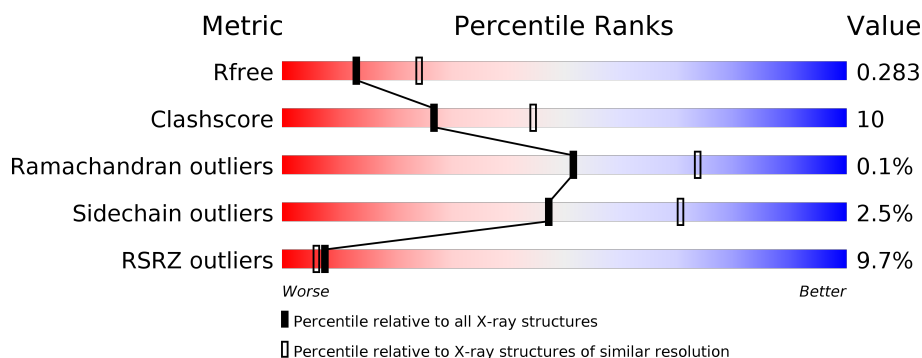
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.57 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	3676 (2.60-2.56)
Clashscore	141614	4049 (2.60-2.56)
Ramachandran outliers	138981	3979 (2.60-2.56)
Sidechain outliers	138945	3979 (2.60-2.56)
RSRZ outliers	127900	3614 (2.60-2.56)

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 on 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	190	
1	C	190	
2	B	218	
2	D	218	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5900 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 Autophagy-related protein 13.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	183	Total	C	N	O	S	0	0	0
			1343	856	230	249	8			
1	C	183	Total	C	N	O	S	0	0	0
			1365	871	230	256	8			

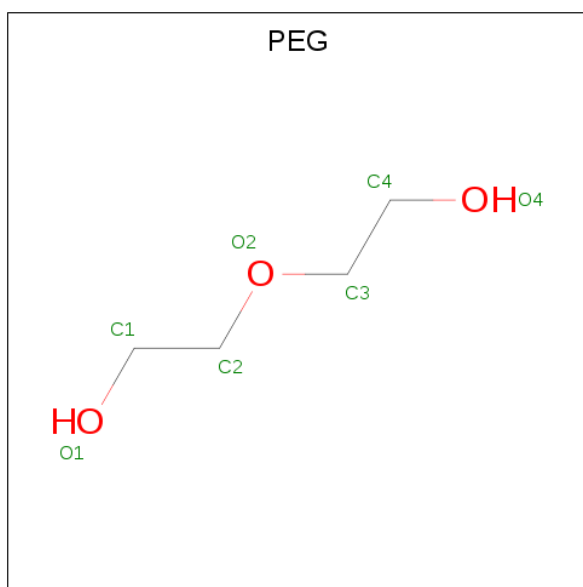
- Molecule 2 is a protein called Autophagy-related protein 101.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	204	Total	C	N	O	S	0	0	0
			1569	995	263	302	9			
2	D	204	Total	C	N	O	S	0	0	0
			1546	983	260	294	9			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	40	ALA	LYS	engineered mutation	UNP Q9BSB4
B	41	ALA	LYS	engineered mutation	UNP Q9BSB4
B	42	ALA	GLU	engineered mutation	UNP Q9BSB4
D	40	ALA	LYS	engineered mutation	UNP Q9BSB4
D	41	ALA	LYS	engineered mutation	UNP Q9BSB4
D	42	ALA	GLU	engineered mutation	UNP Q9BSB4

- Molecule 3 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).



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

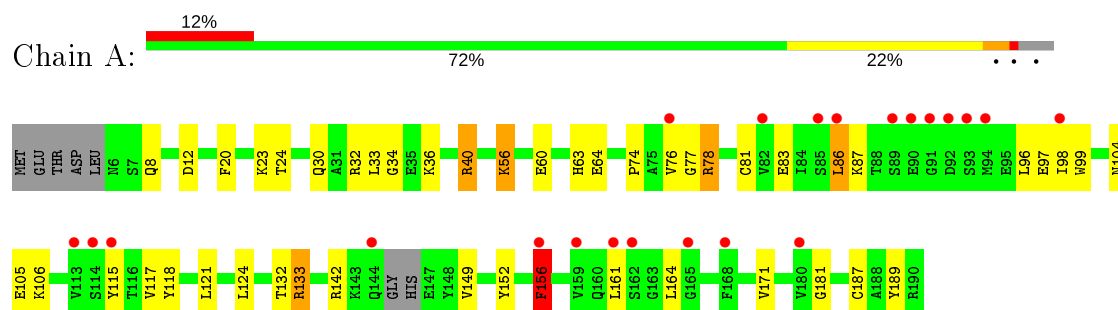
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	12	Total	O	0	0
			12	12		
4	B	15	Total	O	0	0
			15	15		
4	C	25	Total	O	0	0
			25	25		
4	D	18	Total	O	0	0
			18	18		

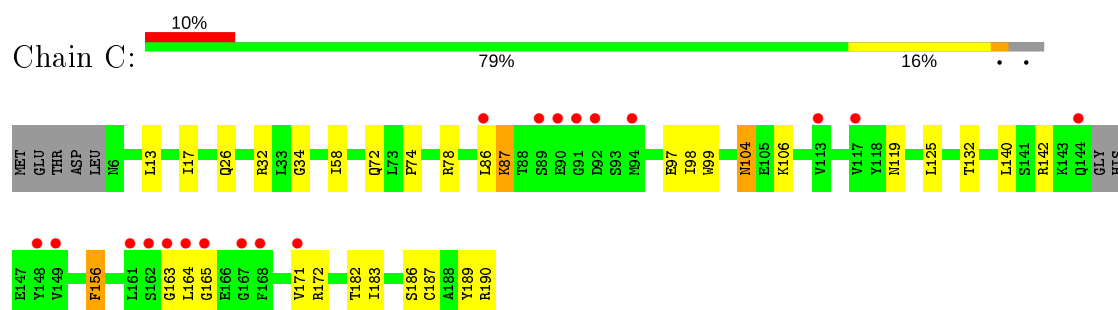
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

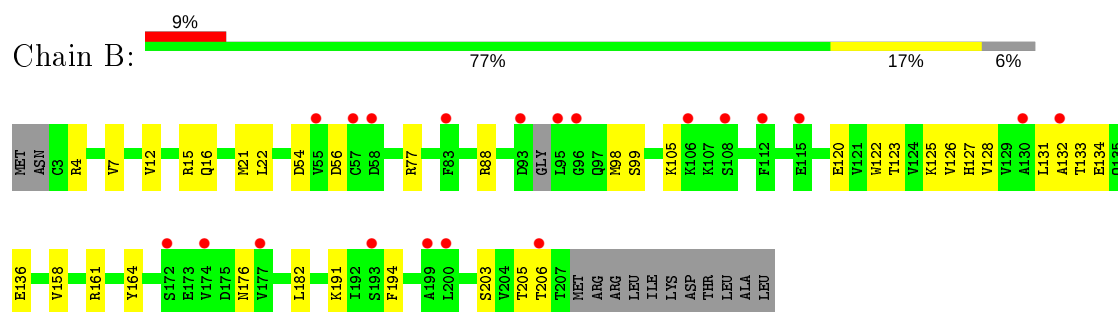
• Molecule 1: Autophagy-related protein 13



• Molecule 1: Autophagy-related protein 13

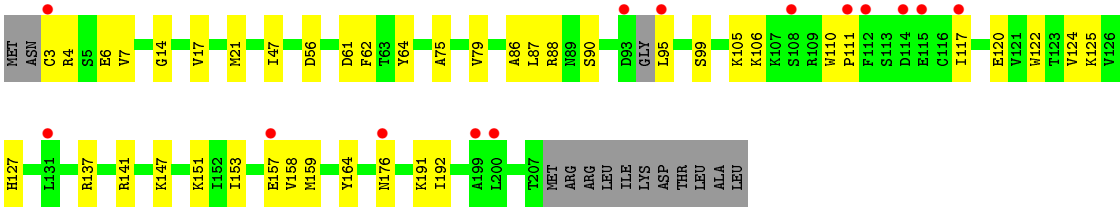


• Molecule 2: Autophagy-related protein 101



• Molecule 2: Autophagy-related protein 101





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 ₁ 2 ₁ 2	Depositor
Cell constants a, b, c, α , β , γ	86.78Å 101.12Å 125.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.59 – 2.57 43.39 – 2.55	Depositor EDS
% Data completeness (in resolution range)	98.8 (38.59-2.57) 87.3 (43.39-2.55)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	0.38 (at 2.54Å)	Xtriage
Refinement program	PHENIX 1.10.1 _2155	Depositor
R, R_{free}	0.232 , 0.283 0.232 , 0.283	Depositor DCC
R_{free} test set	2000 reflections (5.51%)	wwPDB-VP
Wilson B-factor (Å ²)	49.3	Xtriage
Anisotropy	0.871	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 58.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.53$, $\langle L^2 \rangle = 0.37$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5900	wwPDB-VP
Average B, all atoms (Å ²)	82.0	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 46.40 % 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 1.1505e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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

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	0.46	4/1363 (0.3%)	0.63	4/1852 (0.2%)
1	C	0.26	0/1386	0.48	0/1880
2	B	0.30	0/1598	0.49	0/2169
2	D	0.26	0/1575	0.52	0/2140
All	All	0.32	4/5922 (0.1%)	0.53	4/8041 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	40	ARG	CZ-NH1	-6.85	1.24	1.33
1	A	40	ARG	NE-CZ	-6.64	1.24	1.33
1	A	40	ARG	CZ-NH2	-6.35	1.24	1.33
1	A	40	ARG	CD-NE	-5.12	1.37	1.46

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	78	ARG	NE-CZ-NH1	-9.37	115.61	120.30
1	A	156	PHE	CB-CG-CD1	6.12	125.09	120.80
1	A	78	ARG	NE-CZ-NH2	5.78	123.19	120.30
1	A	78	ARG	CG-CD-NE	5.20	122.71	111.80

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	1343	0	1301	31	0
1	C	1365	0	1341	24	0
2	B	1569	0	1490	32	0
2	D	1546	0	1457	33	0
3	C	7	0	10	3	0
4	A	12	0	0	1	0
4	B	15	0	0	3	0
4	C	25	0	0	0	0
4	D	18	0	0	1	0
All	All	5900	0	5599	116	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 10.

All (116) 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:A:32:ARG:NH2	1:A:132:THR:O	2.03	0.91
1:C:104:ASN:ND2	1:C:182:THR:OG1	2.11	0.82
2:B:120:GLU:OE1	2:B:122:TRP:NE1	2.14	0.81
1:C:32:ARG:NH2	1:C:132:THR:O	2.16	0.79
1:C:142:ARG:NH2	2:D:56:ASP:OD1	2.18	0.76
1:A:74:PRO:HA	1:A:156:PHE:HZ	1.52	0.73
2:D:79:VAL:HG11	2:D:192:ILE:HD11	1.70	0.73
1:A:60:GLU:OE2	1:A:152:TYR:OH	2.06	0.72
2:B:4:ARG:NH2	2:B:164:TYR:O	2.23	0.71
2:D:4:ARG:NH2	2:D:120:GLU:OE1	2.21	0.70
1:A:115:TYR:HA	1:A:118:TYR:HB3	1.74	0.69
2:B:132:ALA:H	2:B:136:GLU:HG3	1.60	0.66
1:A:133:ARG:NH1	2:B:54:ASP:OD2	2.30	0.64
1:A:104:ASN:OD1	1:A:106:LYS:N	2.31	0.64
1:C:72:GLN:HB3	1:C:78:ARG:HH21	1.63	0.62
2:B:12:VAL:CG1	2:B:16:GLN:HB2	2.32	0.60
2:B:7:VAL:HG21	2:B:125:LYS:HE3	1.83	0.60
2:B:133:THR:HG23	2:B:136:GLU:HG2	1.84	0.60
2:D:4:ARG:NH1	2:D:164:TYR:O	2.32	0.60
2:D:61:ASP:OD1	2:D:141:ARG:NH2	2.34	0.60
2:B:131:LEU:HB3	2:B:136:GLU:HB3	1.84	0.60
3:C:201:PEG:H21	2:D:47:ILE:HD12	1.85	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:119:ASN:HD21	3:C:201:PEG:H22	1.68	0.57
1:C:32:ARG:NH1	1:C:97:GLU:OE1	2.37	0.57
1:C:26:GLN:HE22	1:C:58:ILE:H	1.51	0.57
2:B:133:THR:H	2:B:136:GLU:HG3	1.69	0.57
1:C:165:GLY:HA3	1:C:190:ARG:HH12	1.70	0.57
2:B:133:THR:CG2	2:B:136:GLU:HG2	2.36	0.56
2:B:133:THR:H	2:B:136:GLU:CG	2.18	0.56
1:A:20:PHE:O	1:A:24:THR:HG23	2.06	0.55
1:C:142:ARG:HH22	2:D:56:ASP:CG	2.08	0.55
2:B:131:LEU:HB3	2:B:136:GLU:CB	2.36	0.54
1:A:64:GLU:OE2	4:A:201:HOH:O	2.19	0.54
1:C:72:GLN:HB3	1:C:78:ARG:NH2	2.23	0.54
1:A:99:TRP:CE3	1:A:187:CYS:HB2	2.43	0.54
2:B:164:TYR:HD1	4:B:303:HOH:O	1.90	0.54
2:B:105:LYS:HB2	2:B:176:ASN:HA	1.91	0.53
2:D:79:VAL:CG1	2:D:192:ILE:HD11	2.37	0.53
2:D:105:LYS:HB2	2:D:176:ASN:HA	1.91	0.53
1:A:86:LEU:HB2	1:A:96:LEU:HD11	1.91	0.52
2:B:4:ARG:NH2	4:B:303:HOH:O	2.42	0.52
1:A:32:ARG:NH1	1:A:97:GLU:OE2	2.43	0.52
1:C:99:TRP:CE3	1:C:187:CYS:HB2	2.44	0.52
2:D:6:GLU:HG3	2:D:158:VAL:HG11	1.91	0.51
2:B:203:SER:O	4:B:301:HOH:O	2.19	0.51
2:D:14:GLY:O	2:D:17:VAL:HG22	2.11	0.51
1:C:163:GLY:O	1:C:190:ARG:NH2	2.44	0.50
2:D:137:ARG:O	2:D:141:ARG:HG3	2.12	0.49
2:D:17:VAL:O	2:D:21:MET:HG2	2.13	0.49
1:A:8:GLN:NE2	1:A:12:ASP:OD1	2.46	0.49
2:B:21:MET:HE1	2:B:126:VAL:HG21	1.95	0.49
1:A:171:VAL:HG23	1:A:189:TYR:CE1	2.47	0.49
1:C:34:GLY:HA2	1:C:142:ARG:HA	1.94	0.49
1:A:76:VAL:O	1:A:78:ARG:N	2.47	0.48
1:C:86:LEU:HD22	1:C:140:LEU:HD21	1.95	0.48
1:A:171:VAL:HG23	1:A:189:TYR:HE1	1.78	0.48
1:A:161:LEU:HB3	1:A:164:LEU:HD12	1.96	0.48
2:B:205:THR:HA	2:B:206:THR:CB	2.44	0.47
2:B:132:ALA:N	2:B:136:GLU:HG3	2.29	0.47
2:B:21:MET:HG2	2:B:98:MET:HE1	1.96	0.47
2:D:192:ILE:HG13	2:D:192:ILE:O	2.14	0.47
2:B:21:MET:HG2	2:B:98:MET:CE	2.45	0.47
2:D:99:SER:O	2:D:192:ILE:HA	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:40:ARG:HA	1:A:40:ARG:HD2	1.42	0.47
1:C:106:LYS:O	1:C:182:THR:HG23	2.15	0.47
2:D:86:ALA:O	2:D:90:SER:OG	2.31	0.46
2:D:106:LYS:HD2	2:D:117:ILE:HD11	1.98	0.46
1:C:98:ILE:HG12	1:C:164:LEU:HD11	1.98	0.46
1:C:172:ARG:HE	1:C:186:SER:HB3	1.80	0.46
1:C:74:PRO:HA	1:C:156:PHE:HZ	1.80	0.46
1:A:83:GLU:HG3	1:A:98:ILE:HG22	1.98	0.45
2:D:110:TRP:CG	2:D:111:PRO:HA	2.51	0.45
1:A:81:CYS:HA	1:A:99:TRP:O	2.16	0.45
3:C:201:PEG:H32	2:D:47:ILE:HD11	1.98	0.45
1:A:30:GLN:HA	1:A:33:LEU:HD22	1.99	0.45
2:B:98:MET:HG3	2:B:194:PHE:HB3	1.99	0.45
1:A:117:VAL:O	1:A:121:LEU:HG	2.16	0.45
1:C:171:VAL:HG13	1:C:189:TYR:HE1	1.81	0.45
1:A:104:ASN:OD1	1:A:105:GLU:N	2.50	0.44
2:D:7:VAL:HG11	2:D:125:LYS:HE2	1.99	0.44
1:C:87:LYS:HG3	1:C:87:LYS:O	2.16	0.44
2:D:88:ARG:NH1	2:D:88:ARG:HB2	2.33	0.44
2:B:12:VAL:O	2:B:128:VAL:HA	2.17	0.44
1:A:34:GLY:HA2	1:A:142:ARG:HA	2.01	0.43
1:A:87:LYS:O	1:A:149:VAL:HG12	2.19	0.43
2:B:133:THR:HG23	2:B:136:GLU:H	1.84	0.43
2:B:133:THR:OG1	2:B:134:GLU:N	2.51	0.43
2:D:191:LYS:NZ	4:D:302:HOH:O	2.52	0.43
1:C:17:ILE:HD12	1:C:183:ILE:HD11	2.01	0.43
2:B:133:THR:N	2:B:136:GLU:HG3	2.33	0.42
2:B:22:LEU:HD11	2:B:77:ARG:HG2	2.01	0.42
1:C:13:LEU:O	1:C:17:ILE:HG12	2.19	0.42
2:D:4:ARG:NH2	2:D:120:GLU:HB2	2.35	0.42
2:B:182:LEU:HD12	2:B:182:LEU:HA	1.57	0.42
2:D:99:SER:HA	2:D:122:TRP:O	2.19	0.42
2:D:159:MET:HB3	2:D:159:MET:HE2	1.84	0.42
1:A:36:LYS:HD2	2:B:56:ASP:OD2	2.19	0.42
1:A:30:GLN:NE2	1:A:56:LYS:O	2.47	0.42
2:D:153:ILE:O	2:D:157:GLU:HG3	2.18	0.42
2:D:88:ARG:HH11	2:D:88:ARG:HB2	1.84	0.42
1:C:171:VAL:HG13	1:C:189:TYR:CE1	2.55	0.42
2:D:3:CYS:N	2:D:4:ARG:HA	2.34	0.42
1:A:20:PHE:CZ	1:A:124:LEU:HD23	2.55	0.41
1:A:33:LEU:HA	1:A:33:LEU:HD12	1.91	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:147:LYS:O	2:D:151:LYS:HG2	2.21	0.41
1:A:74:PRO:HA	1:A:156:PHE:CZ	2.41	0.41
2:B:158:VAL:HG12	2:B:161:ARG:NH1	2.35	0.41
2:D:62:PHE:HE1	2:D:64:TYR:HB3	1.86	0.41
2:D:14:GLY:HA3	2:D:88:ARG:HH22	1.86	0.41
2:D:87:LEU:CD2	2:D:95:LEU:HB2	2.51	0.41
1:A:106:LYS:O	1:A:181:GLY:HA2	2.20	0.41
2:B:7:VAL:HG23	2:B:123:THR:HB	2.03	0.41
1:A:98:ILE:O	1:A:98:ILE:HG13	2.21	0.40
2:B:127:HIS:CD2	2:B:127:HIS:C	2.95	0.40
1:C:125:LEU:HA	1:C:125:LEU:HD23	1.83	0.40
2:D:75:ALA:O	2:D:79:VAL:HG12	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	179/190 (94%)	172 (96%)	6 (3%)	1 (1%)	25	45
1	C	179/190 (94%)	173 (97%)	6 (3%)	0	100	100
2	B	200/218 (92%)	192 (96%)	8 (4%)	0	100	100
2	D	200/218 (92%)	196 (98%)	4 (2%)	0	100	100
All	All	758/816 (93%)	733 (97%)	24 (3%)	1 (0%)	51	73

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	77	GLY

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	135/167 (81%)	129 (96%)	6 (4%)	28	51
1	C	142/167 (85%)	139 (98%)	3 (2%)	53	75
2	B	165/194 (85%)	161 (98%)	4 (2%)	49	72
2	D	159/194 (82%)	157 (99%)	2 (1%)	69	85
All	All	601/722 (83%)	586 (98%)	15 (2%)	47	70

All (15) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	LYS
1	A	56	LYS
1	A	63	HIS
1	A	86	LEU
1	A	133	ARG
1	A	156	PHE
2	B	15	ARG
2	B	88	ARG
2	B	99	SER
2	B	191	LYS
1	C	87	LYS
1	C	104	ASN
1	C	156	PHE
2	D	124	VAL
2	D	127	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	127	HIS
1	C	26	GLN
1	C	104	ASN
1	C	119	ASN

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Mol	Chain	Res	Type
1	C	144	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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	$\# Z > 2$
3	PEG	C	201	-	6,6,6	0.48	0	5,5,5	0.27	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
3	PEG	C	201	-	-	3/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (3) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	C	201	PEG	O1-C1-C2-O2
3	C	201	PEG	C1-C2-O2-C3
3	C	201	PEG	C4-C3-O2-C2

There are no ring outliers.

1 monomer is involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	201	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 ⓘ

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	183/190 (96%)	0.73	22 (12%) 4 3	50, 85, 146, 191	0
1	C	183/190 (96%)	0.57	19 (10%) 6 5	40, 74, 127, 156	0
2	B	204/218 (93%)	0.77	20 (9%) 7 6	45, 83, 128, 162	0
2	D	204/218 (93%)	0.55	14 (6%) 16 14	41, 78, 133, 161	0
All	All	774/816 (94%)	0.66	75 (9%) 7 6	40, 80, 134, 191	0

All (75) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	112	PHE	5.5
2	D	115	GLU	5.2
1	A	92	ASP	5.1
1	A	159	VAL	5.0
2	D	112	PHE	4.9
1	A	90	GLU	4.6
1	A	114	SER	4.6
1	A	113	VAL	4.6
1	C	162	SER	4.5
2	B	93	ASP	4.4
1	C	168	PHE	4.1
1	C	86	LEU	4.0
1	C	90	GLU	3.9
1	A	94	MET	3.9
1	A	82	VAL	3.8
2	B	132	ALA	3.7
1	C	148	TYR	3.6
1	A	93	SER	3.5
2	B	199	ALA	3.4
1	C	89	SER	3.4
2	B	106	LYS	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	156	PHE	3.3
2	D	3	CYS	3.2
2	B	130	ALA	3.2
1	A	161	LEU	3.1
1	A	86	LEU	3.1
1	A	98	ILE	3.1
2	B	177	VAL	3.1
1	C	167	GLY	3.0
1	C	164	LEU	3.0
2	B	96	GLY	3.0
2	B	172	SER	2.9
1	C	144	GLN	2.9
2	B	95	LEU	2.9
2	D	95	LEU	2.8
1	A	85	SER	2.8
1	A	162	SER	2.8
2	D	131	LEU	2.7
2	B	58	ASP	2.7
1	C	163	GLY	2.7
2	B	115	GLU	2.6
2	B	83	PHE	2.6
2	D	117	ILE	2.6
1	C	165	GLY	2.6
1	A	180	VAL	2.6
2	B	174	VAL	2.6
1	C	149	VAL	2.5
1	A	89	SER	2.5
2	D	111	PRO	2.5
2	B	193	SER	2.4
1	A	144	GLN	2.4
1	C	113	VAL	2.4
2	D	108	SER	2.4
2	B	108	SER	2.4
1	A	91	GLY	2.3
1	C	171	VAL	2.3
2	B	57	CYS	2.3
2	D	199	ALA	2.3
1	C	161	LEU	2.3
1	A	168	PHE	2.2
2	B	200	LEU	2.2
1	A	115	TYR	2.2
2	D	157	GLU	2.2

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Mol	Chain	Res	Type	RSRZ
2	B	55	VAL	2.2
2	D	93	ASP	2.1
2	D	176	ASN	2.1
2	D	114	ASP	2.1
1	A	165	GLY	2.1
1	C	94	MET	2.1
1	A	76	VAL	2.0
2	D	200	LEU	2.0
1	C	117	VAL	2.0
1	C	91	GLY	2.0
2	B	206	THR	2.0
1	C	92	ASP	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 carbohydrates 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
3	PEG	C	201	7/7	0.81	0.21	71,85,91,92	0

6.5 Other polymers [i](#)

There are no such residues in this entry.