



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 08:01 PM UTC

PDB ID : 3C1L / pdb_00003c1l
Title : Crystal structure of an antioxidant defense protein (mlr4105) from mesorhizobium loti maff303099 at 2.00 Å resolution
Authors : Joint Center for Structural Genomics (JCSG)
Deposited on : 2008-01-23
Resolution : 2.00 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

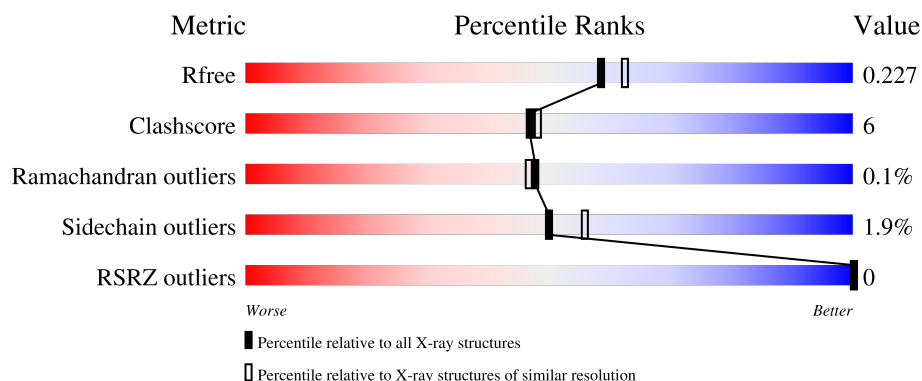
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	10052 (2.00-2.00)
Clashscore	190562	11152 (2.00-2.00)
Ramachandran outliers	187476	11031 (2.00-2.00)
Sidechain outliers	187428	11029 (2.00-2.00)
RSRZ outliers	180081	10067 (2.00-2.00)

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	188	
1	B	188	
1	C	188	
1	D	188	
1	E	188	

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Mol	Chain	Length	Quality of chain
1	F	188	 88% 7% . .
1	G	188	 87% 11% . .
1	H	188	 86% 8% . .
1	I	188	 87% 9% . .
1	J	188	 84% 13% . .
1	K	188	 81% 14% . .
1	L	188	 81% 14% . .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PEG	I	190	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 17569 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 Putative antioxidant defense protein Mlr4105.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	182	Total	C	N	O	S	Se	0	3	0
			1408	887	243	267	3	8			
1	B	185	Total	C	N	O	S	Se	0	1	0
			1423	894	245	273	3	8			
1	C	183	Total	C	N	O	S	Se	0	3	0
			1414	891	246	266	3	8			
1	D	180	Total	C	N	O	S	Se	0	1	0
			1391	876	241	263	3	8			
1	E	181	Total	C	N	O	S	Se	0	0	0
			1392	876	239	266	3	8			
1	F	181	Total	C	N	O	S	Se	0	0	0
			1383	870	238	264	3	8			
1	G	185	Total	C	N	O	S	Se	0	0	0
			1416	893	244	267	3	9			
1	H	180	Total	C	N	O	S	Se	0	0	0
			1359	856	233	259	3	8			
1	I	181	Total	C	N	O	S	Se	0	1	0
			1393	877	241	264	3	8			
1	J	183	Total	C	N	O	S	Se	0	0	0
			1400	882	241	266	3	8			
1	K	181	Total	C	N	O	S	Se	0	0	0
			1354	848	237	258	3	8			
1	L	181	Total	C	N	O	S	Se	0	0	0
			1329	833	231	254	3	8			

There are 12 discrepancies between the modelled and reference sequences:

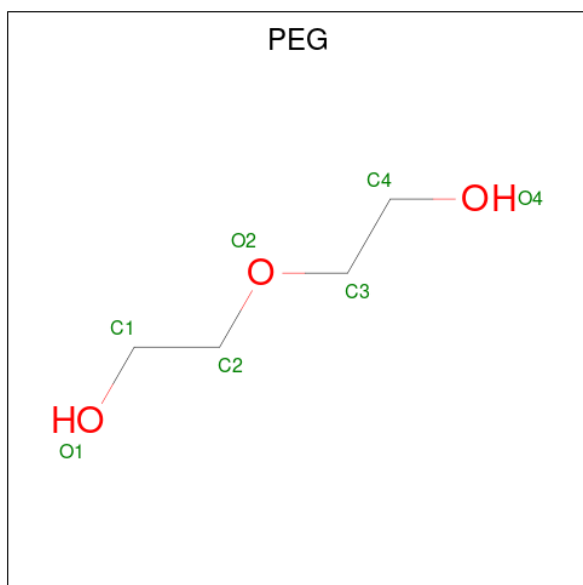
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q98ES4
B	0	GLY	-	expression tag	UNP Q98ES4
C	0	GLY	-	expression tag	UNP Q98ES4
D	0	GLY	-	expression tag	UNP Q98ES4
E	0	GLY	-	expression tag	UNP Q98ES4

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Chain	Residue	Modelled	Actual	Comment	Reference
F	0	GLY	-	expression tag	UNP Q98ES4
G	0	GLY	-	expression tag	UNP Q98ES4
H	0	GLY	-	expression tag	UNP Q98ES4
I	0	GLY	-	expression tag	UNP Q98ES4
J	0	GLY	-	expression tag	UNP Q98ES4
K	0	GLY	-	expression tag	UNP Q98ES4
L	0	GLY	-	expression tag	UNP Q98ES4

- Molecule 2 is DI(HYDROXYETHYL)ETHER (CCD ID: PEG) (formula: C₄H₁₀O₃).



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

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	C	1	Total C O 7 4 3	0	0
2	C	1	Total C O 7 4 3	0	0
2	C	1	Total C O 7 4 3	0	0
2	D	1	Total C O 7 4 3	0	0
2	D	1	Total C O 7 4 3	0	0
2	D	1	Total C O 7 4 3	0	0
2	E	1	Total C O 7 4 3	0	0
2	E	1	Total C O 7 4 3	0	0
2	G	1	Total C O 7 4 3	0	0
2	H	1	Total C O 7 4 3	0	0
2	H	1	Total C O 7 4 3	0	0
2	I	1	Total C O 7 4 3	0	0
2	I	1	Total C O 7 4 3	0	0
2	I	1	Total C O 7 4 3	0	0
2	J	1	Total C O 7 4 3	0	0
2	J	1	Total C O 7 4 3	0	0
2	J	1	Total C O 7 4 3	0	0

- Molecule 3 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	J	1	Total Cl 1 1	0	0

- Molecule 4 is water.

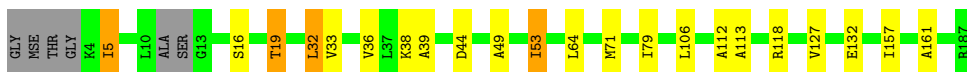
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	75	Total 75	O 75	0	0
4	B	76	Total 76	O 76	0	0
4	C	76	Total 76	O 76	0	0
4	D	87	Total 87	O 87	0	0
4	E	78	Total 78	O 78	0	0
4	F	64	Total 64	O 64	0	0
4	G	57	Total 57	O 57	0	0
4	H	48	Total 48	O 48	0	0
4	I	56	Total 56	O 56	0	0
4	J	59	Total 59	O 59	0	0
4	K	33	Total 33	O 33	0	0
4	L	22	Total 22	O 22	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: Putative antioxidant defense protein Mlr4105

Chain A: 



- Molecule 1: Putative antioxidant defense protein Mlr4105

Chain B: 



- Molecule 1: Putative antioxidant defense protein Mlr4105

Chain C: 



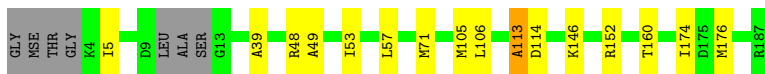
- Molecule 1: Putative antioxidant defense protein Mlr4105

Chain D: 



- Molecule 1: Putative antioxidant defense protein Mlr4105

Chain E: 



- Molecule 1: Putative antioxidant defense protein Mlr4105

Chain F: 



- Molecule 1: Putative antioxidant defense protein Mlr4105

Chain G: 87% 11% ..



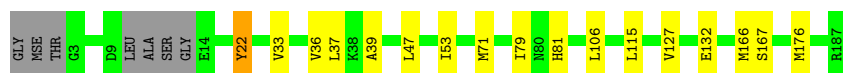
- Molecule 1: Putative antioxidant defense protein Mlr4105

Chain H: 86% 8% ..



- Molecule 1: Putative antioxidant defense protein Mlr4105

Chain I: 87% 9% ..



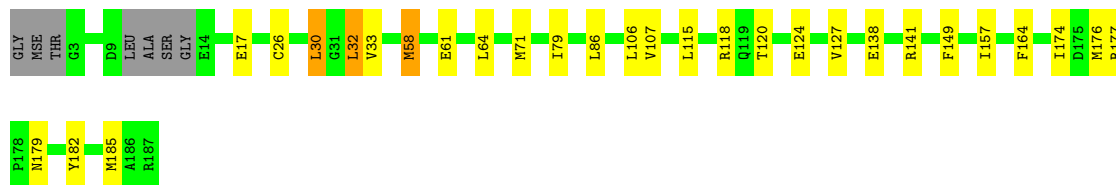
- Molecule 1: Putative antioxidant defense protein Mlr4105

Chain J: 84% 13% ..



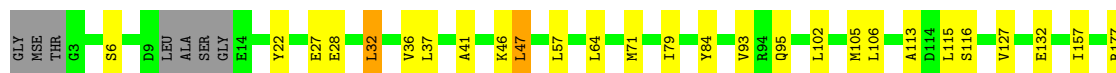
- Molecule 1: Putative antioxidant defense protein Mlr4105

Chain K: 81% 14% ..



- Molecule 1: Putative antioxidant defense protein Mlr4105

Chain L: 81% 14% ..



M185
A186
R187

4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	65.69Å 68.06Å 266.83Å 90.00° 95.94° 90.00°	Depositor
Resolution (Å)	29.81 – 2.00 29.81 – 2.00	Depositor EDS
% Data completeness (in resolution range)	87.5 (29.81-2.00) 87.5 (29.81-2.00)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.84 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.4.0066, PHENIX	Depositor
R, R_{free}	0.171 , 0.225 0.176 , 0.227	Depositor DCC
R_{free} test set	6959 reflections (4.39%)	wwPDB-VP
Wilson B-factor (Å ²)	25.3	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 39.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.045 for h,-k,-h-l	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	17569	wwPDB-VP
Average B, all atoms (Å ²)	30.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.89% 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: CL, 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	1.08	1/1432 (0.1%)	1.10	1/1920 (0.1%)
1	B	1.04	0/1439	1.09	4/1928 (0.2%)
1	C	1.10	1/1438 (0.1%)	1.09	1/1927 (0.1%)
1	D	1.06	0/1409	1.05	1/1886 (0.1%)
1	E	1.03	0/1407	1.02	1/1884 (0.1%)
1	F	0.98	0/1398	1.00	1/1874 (0.1%)
1	G	1.01	0/1431	1.06	0/1915
1	H	0.89	0/1374	1.02	2/1845 (0.1%)
1	I	1.02	2/1411 (0.1%)	1.01	0/1890
1	J	0.94	1/1415 (0.1%)	1.04	2/1895 (0.1%)
1	K	0.75	0/1369	0.93	0/1837
1	L	0.87	2/1343 (0.1%)	0.96	0/1808
All	All	0.99	7/16866 (0.0%)	1.03	13/22609 (0.1%)

All (7) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	L	28	GLU	CD-OE2	10.06	1.44	1.25
1	L	28	GLU	CD-OE1	7.25	1.39	1.25
1	I	22	TYR	N-CA	-5.93	1.39	1.46
1	A	161	ALA	CA-CB	-5.59	1.44	1.53
1	J	157	ILE	CA-CB	5.51	1.61	1.54
1	C	87	THR	CA-CB	5.24	1.61	1.53
1	I	33	VAL	CA-CB	5.15	1.59	1.53

All (13) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	J	30	LEU	N-CA-C	-8.30	98.55	110.59
1	C	53	ILE	CB-CA-C	-5.57	104.72	112.02

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	53	ILE	CB-CA-C	-5.51	104.61	112.22
1	H	53	ILE	CB-CA-C	-5.48	104.66	112.22
1	D	113	ALA	N-CA-C	5.39	119.57	112.89
1	J	131	GLU	N-CA-C	5.36	117.20	111.36
1	F	30	LEU	N-CA-C	-5.34	102.85	110.59
1	A	113	ALA	N-CA-C	5.31	119.61	113.18
1	H	113	ALA	N-CA-C	5.16	119.58	113.28
1	E	113	ALA	N-CA-C	5.16	117.81	111.82
1	B	25	LYS	N-CA-C	5.05	116.87	111.36
1	B	116	SER	CA-C-N	5.02	125.05	119.32
1	B	116	SER	C-N-CA	5.02	125.05	119.32

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	1408	0	1392	16	0
1	B	1423	0	1396	18	0
1	C	1414	0	1396	16	0
1	D	1391	0	1376	17	0
1	E	1392	0	1369	22	0
1	F	1383	0	1352	19	0
1	G	1416	0	1407	16	0
1	H	1359	0	1312	15	0
1	I	1393	0	1369	15	0
1	J	1400	0	1379	18	0
1	K	1354	0	1288	21	0
1	L	1329	0	1243	24	0
2	A	28	0	40	1	0
2	B	28	0	40	4	0
2	C	21	0	30	1	0
2	D	21	0	30	1	0
2	E	14	0	20	0	0
2	G	7	0	10	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	H	14	0	20	0	0
2	I	21	0	30	5	0
2	J	21	0	30	0	0
3	J	1	0	0	0	0
4	A	75	0	0	1	0
4	B	76	0	0	0	0
4	C	76	0	0	1	0
4	D	87	0	0	1	0
4	E	78	0	0	1	0
4	F	64	0	0	0	0
4	G	57	0	0	2	0
4	H	48	0	0	1	0
4	I	56	0	0	1	0
4	J	59	0	0	0	0
4	K	33	0	0	2	0
4	L	22	0	0	2	0
All	All	17569	0	16529	189	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 6.

All (189) 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:E:53:ILE:HD12	1:F:53:ILE:CD1	1.48	1.44
1:E:53:ILE:CD1	1:F:53:ILE:CD1	2.25	1.15
1:E:53:ILE:CD1	1:F:53:ILE:HD12	1.84	1.06
1:E:53:ILE:HD12	1:F:53:ILE:HD11	1.05	1.02
1:G:53:ILE:HG13	1:H:53:ILE:HD11	1.42	1.02
1:E:53:ILE:HD12	1:F:53:ILE:HD12	1.42	0.93
1:A:16:SER:OG	1:A:19[A]:THR:HG23	1.70	0.91
1:K:79:ILE:HD11	1:K:127:VAL:HG22	1.59	0.84
1:B:71:MSE:HG2	1:B:106:LEU:HD21	1.61	0.80
1:I:53:ILE:HG13	1:J:53:ILE:HD11	1.65	0.79
1:E:53:ILE:CD1	1:F:53:ILE:HD11	2.00	0.78
1:D:79:ILE:HD11	1:D:127:VAL:HG22	1.66	0.77
1:L:71:MSE:HE3	1:L:106:LEU:HD11	1.68	0.76
1:E:57:LEU:HD21	1:E:160:THR:HA	1.68	0.74
1:J:101:ALA:HB1	1:J:105:MSE:HE3	1.70	0.74
1:E:71:MSE:HE3	1:E:106:LEU:HD21	1.69	0.74
1:G:53:ILE:CG1	1:H:53:ILE:HD11	2.17	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:5:ILE:HD12	1:F:5:ILE:C	2.17	0.70
1:L:71:MSE:HE1	1:L:115:LEU:HD21	1.73	0.69
1:G:79:ILE:HD11	1:G:127:VAL:HG22	1.74	0.69
1:H:79:ILE:HD11	1:H:127:VAL:HG22	1.72	0.69
1:L:93:VAL:HG11	1:L:106:LEU:HD12	1.75	0.69
1:G:114:ASP:HB3	4:G:528:HOH:O	1.93	0.68
1:I:71:MSE:HE1	1:I:115:LEU:HD22	1.77	0.67
1:K:71:MSE:HE1	1:K:115:LEU:HD22	1.76	0.67
1:C:71:MSE:HG2	1:C:106[A]:LEU:CD1	2.26	0.66
1:C:71:MSE:HG2	1:C:106[A]:LEU:HD13	1.79	0.63
1:G:71:MSE:HE2	1:G:106:LEU:HD11	1.80	0.62
1:K:174:ILE:HD11	1:K:176:MSE:HE3	1.81	0.62
1:G:71:MSE:CE	1:G:106:LEU:HD11	2.28	0.62
1:A:118:ARG:NH1	4:A:232:HOH:O	2.32	0.62
1:B:119:GLN:HE22	2:B:191:PEG:C3	2.13	0.62
4:K:711:HOH:O	1:L:46:LYS:CB	2.49	0.61
1:C:36:VAL:HG21	2:C:188:PEG:H32	1.81	0.61
1:J:5:ILE:CD1	1:J:39:ALA:HB2	2.31	0.61
1:A:112:ALA:O	1:C:177[B]:ARG:NH2	2.33	0.60
1:C:53:ILE:HD11	1:D:53:ILE:HG13	1.82	0.60
1:G:15:LEU:HD22	1:G:41:ALA:HB1	1.83	0.60
1:F:47:LEU:O	1:F:51:THR:HG23	2.01	0.59
1:D:19:THR:HG21	1:D:44:ASP:OD1	2.02	0.58
1:B:79:ILE:HD11	1:B:127:VAL:HG22	1.86	0.58
1:F:27:GLU:O	1:F:30:LEU:O	2.22	0.58
1:H:71:MSE:HG2	1:H:106:LEU:HD21	1.85	0.58
1:K:118:ARG:HB2	4:K:717:HOH:O	2.04	0.58
1:L:32:LEU:C	1:L:32:LEU:HD22	2.29	0.58
1:F:5:ILE:HD12	1:F:6:SER:N	2.20	0.57
1:I:36:VAL:HG21	2:I:190:PEG:H21	1.86	0.57
1:C:124:GLU:HA	1:C:124:GLU:OE1	2.05	0.56
1:D:36:VAL:HG21	2:D:188:PEG:H22	1.87	0.56
1:B:71:MSE:HG2	1:B:106:LEU:CD2	2.31	0.56
1:C:93:VAL:HG11	1:C:106[A]:LEU:HD12	1.88	0.56
1:C:57:LEU:HD23	1:C:58:MSE:SE	2.56	0.56
1:A:79:ILE:HD11	1:A:127:VAL:HG22	1.88	0.56
1:J:47:LEU:O	1:J:51:THR:HG23	2.07	0.55
1:K:118:ARG:HG2	1:K:149:PHE:CZ	2.42	0.54
1:L:32:LEU:HD12	1:L:185:MSE:HE1	1.90	0.54
1:H:71:MSE:HG2	1:H:106:LEU:CD2	2.37	0.54
1:A:33:VAL:HB	1:A:38:LYS:HE3	1.89	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:32:LEU:HD12	1:D:32:LEU:C	2.33	0.53
1:B:35:ASN:OD1	1:B:38:LYS:NZ	2.36	0.53
1:K:32:LEU:HD23	1:K:179:ASN:ND2	2.24	0.52
1:L:41:ALA:HA	1:L:47:LEU:HD12	1.90	0.52
1:I:71:MSE:HG2	1:I:106:LEU:CD1	2.40	0.52
1:J:5:ILE:HD13	1:J:39:ALA:HB2	1.91	0.52
1:L:71:MSE:HE2	1:L:106:LEU:HD21	1.92	0.52
1:B:16:SER:OG	1:B:44:ASP:OD2	2.17	0.52
1:K:71:MSE:HE1	1:K:115:LEU:CD2	2.40	0.52
1:K:182:TYR:HA	1:K:185:MSE:HE3	1.91	0.52
1:I:81:HIS:ND1	4:I:624:HOH:O	2.33	0.52
1:L:36:VAL:HG13	1:L:37:LEU:N	2.25	0.51
1:E:49:ALA:HB1	1:F:53:ILE:HG12	1.93	0.51
1:L:116:SER:HB2	4:L:741:HOH:O	2.09	0.51
1:L:71:MSE:HE1	1:L:115:LEU:CD2	2.41	0.51
1:B:53:ILE:O	1:B:57:LEU:HB2	2.11	0.51
1:E:71:MSE:CE	1:E:106:LEU:HD21	2.40	0.51
1:J:123:LEU:O	1:J:127:VAL:HG23	2.11	0.51
1:L:22:TYR:HE2	1:L:37:LEU:HD13	1.74	0.51
1:H:141:ARG:NH2	4:H:554:HOH:O	2.44	0.50
1:I:36:VAL:CG2	2:I:190:PEG:H32	2.41	0.50
1:F:5:ILE:C	1:F:5:ILE:CD1	2.84	0.50
1:D:57:LEU:HD21	1:D:160:THR:HA	1.94	0.49
1:F:71:MSE:HG2	1:F:106:LEU:CD2	2.42	0.49
1:H:17:GLU:HB2	1:H:18:PRO:HD3	1.94	0.49
1:L:71:MSE:HE3	1:L:102:LEU:HD11	1.94	0.49
1:J:112:ALA:HB1	1:K:177:ARG:HH12	1.78	0.49
1:A:19[A]:THR:HG21	1:A:44:ASP:OD1	2.12	0.49
1:C:70:GLU:O	1:C:74:VAL:HG23	2.13	0.49
1:A:36:VAL:HG11	2:A:190:PEG:H31	1.95	0.49
1:F:79:ILE:HD11	1:F:127:VAL:HG22	1.95	0.49
1:L:79:ILE:HD11	1:L:127:VAL:HG22	1.95	0.48
1:C:118[B]:ARG:NH1	1:C:118[B]:ARG:HB3	2.27	0.48
1:L:84:TYR:CD1	1:L:84:TYR:C	2.91	0.48
1:E:113:ALA:O	1:E:114:ASP:HB2	2.14	0.48
1:G:70:GLU:HG3	1:G:96:LEU:HD12	1.96	0.48
1:D:71:MSE:HE1	1:D:115:LEU:HD22	1.94	0.48
1:L:93:VAL:HG11	1:L:106:LEU:CD1	2.43	0.47
1:A:71:MSE:HG2	1:A:106:LEU:HD21	1.97	0.47
1:B:119:GLN:HE22	2:B:191:PEG:H32	1.79	0.47
1:D:48:ARG:O	1:D:49:ALA:C	2.57	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:53:ILE:CD1	1:H:53:ILE:HD11	2.45	0.47
1:J:70:GLU:OE1	1:J:89:HIS:ND1	2.36	0.47
1:L:102:LEU:HD11	1:L:106:LEU:HD11	1.97	0.47
1:E:5:ILE:CD1	1:E:39:ALA:HB2	2.45	0.47
1:J:71:MSE:HG2	1:J:106:LEU:CD1	2.44	0.47
1:K:30:LEU:CD1	1:K:30:LEU:N	2.78	0.47
1:J:105:MSE:SE	1:J:113:ALA:HA	2.65	0.46
1:L:71:MSE:HG2	1:L:106:LEU:HD13	1.97	0.46
1:J:5:ILE:HD11	1:J:39:ALA:HA	1.97	0.46
1:G:59:LEU:HD23	1:G:66:LYS:HD2	1.96	0.46
1:K:138:GLU:HA	1:K:141:ARG:NH1	2.31	0.46
1:B:64:LEU:HD11	1:B:157:ILE:HG13	1.98	0.46
1:K:26:CYS:HB2	1:K:33:VAL:HG22	1.97	0.45
1:C:146:LYS:HD2	4:C:264:HOH:O	2.16	0.45
1:D:71:MSE:HE3	1:D:106:LEU:HD11	1.97	0.45
1:K:58:MSE:SE	1:K:164:PHE:CE2	3.20	0.45
1:K:174:ILE:HD11	1:K:176:MSE:CE	2.44	0.45
1:L:177:ARG:HD2	4:L:733:HOH:O	2.16	0.45
1:G:111:ARG:HG2	4:G:530:HOH:O	2.16	0.44
1:H:32:LEU:HD12	1:H:32:LEU:C	2.43	0.44
1:E:71:MSE:HG2	1:E:106:LEU:CD2	2.48	0.44
1:H:105:MSE:SE	1:H:113:ALA:HA	2.67	0.44
1:G:1:MSE:HE1	1:G:6:SER:C	2.43	0.44
1:E:53:ILE:HG12	1:F:49:ALA:HB1	1.99	0.44
1:I:22:TYR:HE2	1:I:37:LEU:HD13	1.83	0.44
1:A:5:ILE:HD11	1:A:39:ALA:HA	1.99	0.43
1:A:49:ALA:HB1	1:B:53:ILE:CD1	2.48	0.43
1:B:84:TYR:CE2	1:B:178:PRO:HA	2.53	0.43
1:D:21:ALA:O	1:D:24:ALA:HB3	2.18	0.43
1:H:104:GLU:O	1:H:108:MSE:HG3	2.18	0.43
1:B:167:SER:HB3	2:B:189:PEG:H32	1.99	0.43
1:D:113:ALA:O	1:D:114:ASP:HB2	2.17	0.43
1:I:36:VAL:HG23	2:I:190:PEG:H32	2.01	0.43
1:K:120:THR:O	1:K:124:GLU:HG2	2.18	0.43
1:A:49:ALA:HB1	1:B:53:ILE:HD13	2.00	0.43
1:E:53:ILE:CD1	1:F:53:ILE:CG1	2.95	0.43
1:L:105:MSE:SE	1:L:113:ALA:HA	2.68	0.43
1:B:119:GLN:HE22	2:B:191:PEG:H31	1.83	0.43
1:H:24:ALA:O	1:H:28:GLU:HG3	2.19	0.43
1:I:79:ILE:HD11	1:I:127:VAL:HG22	2.01	0.43
1:A:53:ILE:CD1	1:B:53:ILE:HG13	2.49	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:32:LEU:HD23	1:D:181:GLU:HB2	2.01	0.43
1:E:105:MSE:SE	1:E:113:ALA:HA	2.69	0.42
1:J:152:ARG:HA	1:J:152:ARG:HD2	1.95	0.42
1:L:64:LEU:HD11	1:L:157:ILE:HG13	2.01	0.42
1:K:86:LEU:HD22	1:K:107:VAL:HG22	2.00	0.42
1:I:53:ILE:HG13	1:J:53:ILE:CD1	2.43	0.42
1:A:53:ILE:HD11	1:B:53:ILE:HG13	2.01	0.42
1:C:79:ILE:HD11	1:C:127:VAL:HG22	2.01	0.42
1:I:37:LEU:HD22	1:I:47:LEU:HD11	2.01	0.42
1:K:71:MSE:HG2	1:K:106:LEU:HD13	2.02	0.42
1:A:71:MSE:HG2	1:A:106:LEU:CD2	2.50	0.42
1:D:110:PHE:C	1:D:110:PHE:CD1	2.97	0.42
1:L:71:MSE:CE	1:L:106:LEU:HD21	2.49	0.42
1:F:63:GLY:HA3	1:F:153:ASP:OD1	2.20	0.42
1:B:155:TRP:CD1	1:B:155:TRP:C	2.98	0.42
1:H:21:ALA:O	1:H:24:ALA:HB3	2.19	0.42
1:E:152:ARG:NH2	4:E:395:HOH:O	2.53	0.42
1:K:64:LEU:HD11	1:K:157:ILE:HG13	2.02	0.41
1:D:90:GLY:HA3	4:D:284:HOH:O	2.20	0.41
1:J:113:ALA:O	1:J:114:ASP:HB2	2.20	0.41
1:A:32:LEU:HD23	1:A:32:LEU:O	2.21	0.41
1:E:71:MSE:HG2	1:E:106:LEU:HD21	2.01	0.41
1:C:73:ALA:HB2	1:C:164:PHE:CE2	2.55	0.41
1:E:174:ILE:HD11	1:E:176:MSE:HB2	2.02	0.41
1:H:19:THR:HG23	1:H:47:LEU:HD21	2.02	0.41
1:I:39:ALA:CB	1:I:176:MSE:HE2	2.51	0.41
1:G:47:LEU:O	1:G:51:THR:HG23	2.20	0.41
1:K:58:MSE:HE3	1:K:58:MSE:HB2	1.87	0.41
1:I:167:SER:HA	2:I:190:PEG:H41	2.02	0.41
1:J:17:GLU:HB2	1:J:18:PRO:HD3	2.03	0.41
1:G:25:LYS:O	1:G:28:GLU:HB3	2.20	0.41
1:K:30:LEU:N	1:K:30:LEU:HD12	2.36	0.41
1:L:95:GLN:HE21	1:L:186:ALA:HB2	1.84	0.41
1:I:36:VAL:HG21	2:I:190:PEG:H32	2.02	0.41
1:J:71:MSE:HG2	1:J:106:LEU:HD11	2.02	0.41
1:K:58:MSE:SE	1:K:164:PHE:HE2	2.53	0.41
1:B:110:PHE:CD1	1:B:110:PHE:C	2.98	0.40
1:L:32:LEU:C	1:L:32:LEU:CD2	2.93	0.40
1:C:53:ILE:HD11	1:D:53:ILE:CD1	2.51	0.40
1:C:53:ILE:CD1	1:D:53:ILE:CD1	2.99	0.40
1:E:53:ILE:HD11	1:F:53:ILE:HG13	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:53:ILE:HG13	1:H:53:ILE:CD1	2.31	0.40
1:G:33:VAL:HB	1:G:38:LYS:HE3	2.02	0.40
1:A:64:LEU:HD11	1:A:157:ILE:HG13	2.02	0.40
1:C:53:ILE:HD11	1:D:53:ILE:CG1	2.47	0.40
1:E:174:ILE:HD11	1:E:176:MSE:CE	2.52	0.40
1:J:79:ILE:HD11	1:J:127:VAL:HG22	2.03	0.40
1:E:53:ILE:HD11	1:F:53:ILE:CD1	2.39	0.40
1:I:166:MSE:HB2	1:J:166:MSE:HB2	2.03	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	181/188 (96%)	180 (99%)	1 (1%)	0	100	100
1	B	184/188 (98%)	182 (99%)	2 (1%)	0	100	100
1	C	182/188 (97%)	178 (98%)	4 (2%)	0	100	100
1	D	177/188 (94%)	175 (99%)	2 (1%)	0	100	100
1	E	177/188 (94%)	175 (99%)	2 (1%)	0	100	100
1	F	177/188 (94%)	175 (99%)	2 (1%)	0	100	100
1	G	181/188 (96%)	178 (98%)	3 (2%)	0	100	100
1	H	176/188 (94%)	172 (98%)	4 (2%)	0	100	100
1	I	178/188 (95%)	177 (99%)	1 (1%)	0	100	100
1	J	179/188 (95%)	174 (97%)	5 (3%)	0	100	100
1	K	177/188 (94%)	169 (96%)	7 (4%)	1 (1%)	21	17
1	L	177/188 (94%)	166 (94%)	10 (6%)	1 (1%)	21	17
All	All	2146/2256 (95%)	2101 (98%)	43 (2%)	2 (0%)	48	46

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	K	61	GLU
1	L	27	GLU

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	144/139 (104%)	138 (96%)	6 (4%)	26	25
1	B	144/139 (104%)	142 (99%)	2 (1%)	59	66
1	C	143/139 (103%)	143 (100%)	0	100	100
1	D	142/139 (102%)	137 (96%)	5 (4%)	32	32
1	E	142/139 (102%)	140 (99%)	2 (1%)	59	66
1	F	140/139 (101%)	139 (99%)	1 (1%)	76	82
1	G	144/139 (104%)	142 (99%)	2 (1%)	59	66
1	H	135/139 (97%)	134 (99%)	1 (1%)	76	82
1	I	141/139 (101%)	140 (99%)	1 (1%)	76	82
1	J	142/139 (102%)	138 (97%)	4 (3%)	38	41
1	K	131/139 (94%)	127 (97%)	4 (3%)	35	37
1	L	125/139 (90%)	120 (96%)	5 (4%)	28	27
All	All	1673/1668 (100%)	1640 (98%)	33 (2%)	50	54

All (33) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	ILE
1	A	19[A]	THR
1	A	19[B]	THR
1	A	32	LEU
1	A	53	ILE
1	A	132	GLU
1	B	30	LEU

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Mol	Chain	Res	Type
1	B	57	LEU
1	D	5	ILE
1	D	6	SER
1	D	53	ILE
1	D	57	LEU
1	D	146	LYS
1	E	48	ARG
1	E	146	LYS
1	F	48	ARG
1	G	27	GLU
1	G	66	LYS
1	H	106	LEU
1	I	132	GLU
1	J	30	LEU
1	J	48	ARG
1	J	57	LEU
1	J	152	ARG
1	K	17	GLU
1	K	30	LEU
1	K	32	LEU
1	K	58	MSE
1	L	6	SER
1	L	32	LEU
1	L	47	LEU
1	L	57	LEU
1	L	132	GLU

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

Mol	Chain	Res	Type
1	B	119	GLN
1	D	95	GLN
1	L	95	GLN

5.3.3 RNA ⓘ

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 26 ligands modelled in this entry, 1 is monoatomic - leaving 25 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	PEG	D	189	-	6,6,6	0.69	0	5,5,5	0.26	0
2	PEG	B	188	-	6,6,6	0.76	0	5,5,5	0.72	0
2	PEG	I	189	-	6,6,6	0.78	0	5,5,5	0.73	0
2	PEG	C	190	-	6,6,6	0.48	0	5,5,5	0.34	0
2	PEG	B	189	-	6,6,6	0.87	0	5,5,5	0.81	0
2	PEG	J	190	-	6,6,6	0.54	0	5,5,5	0.27	0
2	PEG	J	189	-	6,6,6	0.68	0	5,5,5	0.66	0
2	PEG	D	188	-	6,6,6	0.68	0	5,5,5	0.82	0
2	PEG	A	188	-	6,6,6	0.97	0	5,5,5	0.81	0
2	PEG	E	188	-	6,6,6	0.65	0	5,5,5	0.57	0
2	PEG	D	190	-	6,6,6	0.47	0	5,5,5	0.39	0
2	PEG	I	190	-	6,6,6	0.72	0	5,5,5	0.67	0
2	PEG	B	191	-	6,6,6	0.51	0	5,5,5	0.30	0
2	PEG	J	191	-	6,6,6	0.48	0	5,5,5	0.65	0
2	PEG	G	188	-	6,6,6	0.52	0	5,5,5	0.54	0
2	PEG	C	188	-	6,6,6	0.79	0	5,5,5	0.74	0
2	PEG	A	191	-	6,6,6	0.46	0	5,5,5	0.24	0
2	PEG	H	189	-	6,6,6	0.51	0	5,5,5	0.65	0
2	PEG	B	190	-	6,6,6	0.37	0	5,5,5	0.71	0
2	PEG	E	189	-	6,6,6	0.52	0	5,5,5	0.44	0
2	PEG	I	188	-	6,6,6	0.90	0	5,5,5	0.94	0
2	PEG	C	189	-	6,6,6	0.67	0	5,5,5	0.74	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PEG	H	188	-	6,6,6	0.73	0	5,5,5	0.89	0
2	PEG	A	190	-	6,6,6	0.74	0	5,5,5	0.74	0
2	PEG	A	189	-	6,6,6	0.48	0	5,5,5	0.25	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
2	PEG	D	189	-	-	2/4/4/4	-
2	PEG	B	188	-	-	4/4/4/4	-
2	PEG	I	189	-	-	3/4/4/4	-
2	PEG	C	190	-	-	2/4/4/4	-
2	PEG	B	189	-	-	4/4/4/4	-
2	PEG	J	190	-	-	2/4/4/4	-
2	PEG	J	189	-	-	4/4/4/4	-
2	PEG	D	188	-	-	3/4/4/4	-
2	PEG	A	188	-	-	2/4/4/4	-
2	PEG	E	188	-	-	2/4/4/4	-
2	PEG	D	190	-	-	3/4/4/4	-
2	PEG	I	190	-	-	2/4/4/4	-
2	PEG	B	191	-	-	4/4/4/4	-
2	PEG	J	191	-	-	2/4/4/4	-
2	PEG	G	188	-	-	2/4/4/4	-
2	PEG	C	188	-	-	2/4/4/4	-
2	PEG	A	191	-	-	1/4/4/4	-
2	PEG	H	189	-	-	3/4/4/4	-
2	PEG	B	190	-	-	2/4/4/4	-
2	PEG	E	189	-	-	1/4/4/4	-
2	PEG	I	188	-	-	2/4/4/4	-
2	PEG	C	189	-	-	2/4/4/4	-
2	PEG	H	188	-	-	3/4/4/4	-
2	PEG	A	190	-	-	3/4/4/4	-
2	PEG	A	189	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (62) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	188	PEG	C1-C2-O2-C3
2	C	189	PEG	O1-C1-C2-O2
2	C	190	PEG	O2-C3-C4-O4
2	H	189	PEG	C4-C3-O2-C2
2	E	188	PEG	O2-C3-C4-O4
2	B	188	PEG	O1-C1-C2-O2
2	B	190	PEG	O1-C1-C2-O2
2	D	190	PEG	O1-C1-C2-O2
2	G	188	PEG	O2-C3-C4-O4
2	H	188	PEG	O1-C1-C2-O2
2	H	189	PEG	O1-C1-C2-O2
2	I	189	PEG	O1-C1-C2-O2
2	J	189	PEG	O1-C1-C2-O2
2	D	189	PEG	O1-C1-C2-O2
2	D	190	PEG	O2-C3-C4-O4
2	I	190	PEG	O2-C3-C4-O4
2	H	188	PEG	C1-C2-O2-C3
2	A	189	PEG	O1-C1-C2-O2
2	A	190	PEG	O1-C1-C2-O2
2	A	191	PEG	O2-C3-C4-O4
2	B	191	PEG	O1-C1-C2-O2
2	C	190	PEG	O1-C1-C2-O2
2	I	189	PEG	O2-C3-C4-O4
2	J	191	PEG	O1-C1-C2-O2
2	A	190	PEG	O2-C3-C4-O4
2	I	188	PEG	O2-C3-C4-O4
2	J	189	PEG	O2-C3-C4-O4
2	I	189	PEG	C4-C3-O2-C2
2	B	189	PEG	O2-C3-C4-O4
2	D	188	PEG	O1-C1-C2-O2
2	A	188	PEG	O2-C3-C4-O4
2	H	188	PEG	O2-C3-C4-O4
2	J	190	PEG	O1-C1-C2-O2
2	A	189	PEG	O2-C3-C4-O4
2	B	188	PEG	O2-C3-C4-O4
2	A	190	PEG	C4-C3-O2-C2
2	D	188	PEG	C1-C2-O2-C3
2	B	188	PEG	C4-C3-O2-C2
2	J	191	PEG	C1-C2-O2-C3

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Mol	Chain	Res	Type	Atoms
2	D	190	PEG	C1-C2-O2-C3
2	J	189	PEG	C1-C2-O2-C3
2	H	189	PEG	O2-C3-C4-O4
2	I	190	PEG	O1-C1-C2-O2
2	J	189	PEG	C4-C3-O2-C2
2	C	189	PEG	C4-C3-O2-C2
2	D	189	PEG	C1-C2-O2-C3
2	E	188	PEG	C4-C3-O2-C2
2	B	188	PEG	C1-C2-O2-C3
2	B	190	PEG	C4-C3-O2-C2
2	B	191	PEG	O2-C3-C4-O4
2	B	191	PEG	C4-C3-O2-C2
2	B	189	PEG	O1-C1-C2-O2
2	B	189	PEG	C1-C2-O2-C3
2	B	189	PEG	C4-C3-O2-C2
2	I	188	PEG	C4-C3-O2-C2
2	G	188	PEG	C4-C3-O2-C2
2	B	191	PEG	C1-C2-O2-C3
2	C	188	PEG	O2-C3-C4-O4
2	J	190	PEG	C1-C2-O2-C3
2	A	188	PEG	C4-C3-O2-C2
2	E	189	PEG	O2-C3-C4-O4
2	D	188	PEG	O2-C3-C4-O4

There are no ring outliers.

6 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	189	PEG	1	0
2	D	188	PEG	1	0
2	I	190	PEG	5	0
2	B	191	PEG	3	0
2	C	188	PEG	1	0
2	A	190	PEG	1	0

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	174/188 (92%)	-1.56	0 100 100	16, 30, 38, 48	3 (1%)
1	B	177/188 (94%)	-1.52	0 100 100	16, 29, 37, 48	1 (0%)
1	C	175/188 (93%)	-1.56	0 100 100	16, 29, 38, 48	3 (1%)
1	D	172/188 (91%)	-1.48	0 100 100	17, 30, 38, 51	1 (0%)
1	E	173/188 (92%)	-1.58	0 100 100	25, 29, 38, 43	0
1	F	173/188 (92%)	-1.52	0 100 100	26, 30, 38, 43	0
1	G	176/188 (93%)	-1.54	0 100 100	21, 29, 40, 51	0
1	H	172/188 (91%)	-1.52	0 100 100	26, 29, 36, 41	0
1	I	173/188 (92%)	-1.57	0 100 100	17, 29, 37, 52	1 (0%)
1	J	175/188 (93%)	-1.54	0 100 100	23, 30, 38, 46	0
1	K	173/188 (92%)	-1.42	0 100 100	25, 29, 36, 44	0
1	L	173/188 (92%)	-1.26	0 100 100	26, 29, 36, 43	0
All	All	2086/2256 (92%)	-1.51	0 100 100	16, 29, 38, 52	9 (0%)

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

6.4 Ligands ⓘ

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	PEG	H	189	7/7	0.97	0.05	41,43,48,51	0
2	PEG	J	191	7/7	0.97	0.05	33,42,46,49	0
2	PEG	B	190	7/7	0.98	0.05	40,41,48,51	0
2	PEG	B	191	7/7	0.98	0.04	43,47,52,57	0
2	PEG	C	188	7/7	0.98	0.05	33,37,41,49	0
2	PEG	C	189	7/7	0.98	0.04	29,39,45,47	0
2	PEG	D	188	7/7	0.98	0.04	33,38,43,43	0
2	PEG	D	190	7/7	0.98	0.04	39,45,49,51	0
2	PEG	A	191	7/7	0.98	0.04	29,32,43,47	0
2	PEG	I	189	7/7	0.98	0.04	40,42,48,51	0
2	PEG	J	189	7/7	0.98	0.04	31,37,47,47	0
2	PEG	B	188	7/7	0.98	0.04	25,37,42,50	0
2	PEG	D	189	7/7	0.99	0.03	34,44,46,46	0
2	PEG	A	188	7/7	0.99	0.04	19,31,37,41	0
2	PEG	E	188	7/7	0.99	0.04	33,37,44,49	0
2	PEG	E	189	7/7	0.99	0.04	52,54,57,58	0
2	PEG	G	188	7/7	0.99	0.04	28,35,39,49	0
2	PEG	H	188	7/7	0.99	0.04	35,46,53,53	0
2	PEG	A	189	7/7	0.99	0.04	36,36,39,48	0
2	PEG	I	188	7/7	0.99	0.04	32,37,42,43	0
2	PEG	B	189	7/7	0.99	0.04	27,35,42,43	0
2	PEG	I	190	7/7	0.99	0.04	30,36,42,44	0
2	PEG	C	190	7/7	0.99	0.03	36,41,48,49	0
2	PEG	J	190	7/7	0.99	0.04	43,48,56,56	0
2	PEG	A	190	7/7	0.99	0.04	32,40,44,49	0
3	CL	J	188	1/1	1.00	0.02	50,50,50,50	0

6.5 Other polymers ⓘ

There are no such residues in this entry.