



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 15, 2017 – 02:45 am GMT

PDB ID : 1F33
Title : THE STRUCTURAL BASIS FOR DNA PROTECTION BY E. COLI DPS PROTEIN
Authors : Luo, J.; Liu, D.; White, M.A.; Fox, R.O.
Deposited on : 2000-05-31
Resolution : 2.60 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

<http://wwpdb.org/validation/2016/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:

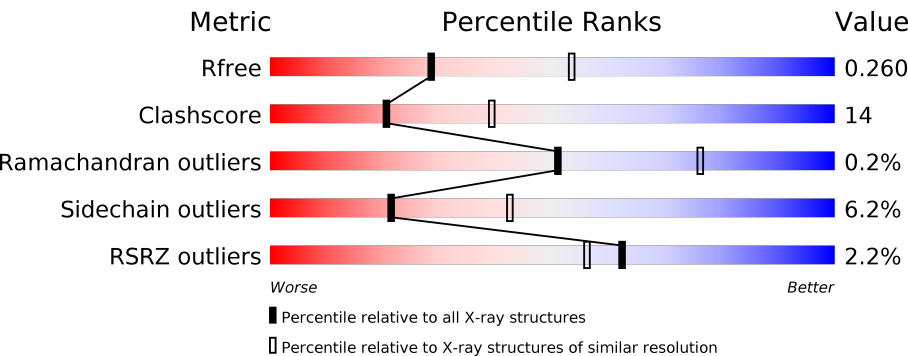
MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

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 2.60 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}	100719	2542 (2.60-2.60)
Clashscore	112137	2895 (2.60-2.60)
Ramachandran outliers	110173	2848 (2.60-2.60)
Sidechain outliers	110143	2848 (2.60-2.60)
RSRZ outliers	101464	2550 (2.60-2.60)

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. 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	167	<div><div>%</div><div><div></div><div>70%</div><div>21%</div><div>••</div><div>7%</div></div></div>
1	B	167	<div><div>%</div><div><div></div><div>68%</div><div>22%</div><div>•</div><div>7%</div></div></div>
1	C	167	<div><div></div><div><div></div><div>68%</div><div>22%</div><div>•</div><div>7%</div></div></div>
1	D	167	<div><div>%</div><div><div></div><div>69%</div><div>22%</div><div>••</div><div>7%</div></div></div>
1	E	167	<div><div>%</div><div><div></div><div>66%</div><div>25%</div><div>••</div><div>7%</div></div></div>
1	F	167	<div><div>%</div><div><div></div><div>66%</div><div>25%</div><div>•</div><div>5%</div></div></div>

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Mol	Chain	Length	Quality of chain
1	G	167	
1	H	167	
1	I	167	
1	J	167	
1	K	167	
1	L	167	

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	TRS	C	603	-	-	-	X
2	TRS	F	606	-	-	-	X
2	TRS	G	607	-	-	-	X

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 15354 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 DNA PROTECTION DURING STARVATION PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	156	Total	C	N	O	S	0	0	0
			1236	776	216	241	3			
1	B	156	Total	C	N	O	S	0	0	0
			1236	776	216	241	3			
1	C	156	Total	C	N	O	S	0	0	0
			1236	776	216	241	3			
1	D	156	Total	C	N	O	S	0	0	0
			1236	776	216	241	3			
1	E	156	Total	C	N	O	S	0	0	0
			1236	776	216	241	3			
1	F	159	Total	C	N	O	S	0	0	0
			1256	788	220	245	3			
1	G	156	Total	C	N	O	S	0	0	0
			1236	776	216	241	3			
1	H	156	Total	C	N	O	S	0	0	0
			1236	776	216	241	3			
1	I	159	Total	C	N	O	S	0	0	0
			1256	788	220	245	3			
1	J	156	Total	C	N	O	S	0	0	0
			1236	776	216	241	3			
1	K	156	Total	C	N	O	S	0	0	0
			1236	776	216	241	3			
1	L	156	Total	C	N	O	S	0	0	0
			1236	776	216	241	3			

- Molecule 2 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: C₄H₁₂NO₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	E	1	Total	C	N	O	0	0
			8	4	1	3		
2	A	1	Total	C	N	O	0	0
			8	4	1	3		
2	C	1	Total	C	N	O	0	0
			8	4	1	3		
2	C	1	Total	C	N	O	0	0
			8	4	1	3		
2	F	1	Total	C	N	O	0	0
			8	4	1	3		
2	F	1	Total	C	N	O	0	0
			8	4	1	3		
2	G	1	Total	C	N	O	0	0
			8	4	1	3		
2	L	1	Total	C	N	O	0	0
			8	4	1	3		
2	I	1	Total	C	N	O	0	0
			8	4	1	3		
2	K	1	Total	C	N	O	0	0
			8	4	1	3		
2	K	1	Total	C	N	O	0	0
			8	4	1	3		
2	J	1	Total	C	N	O	0	0
			8	4	1	3		

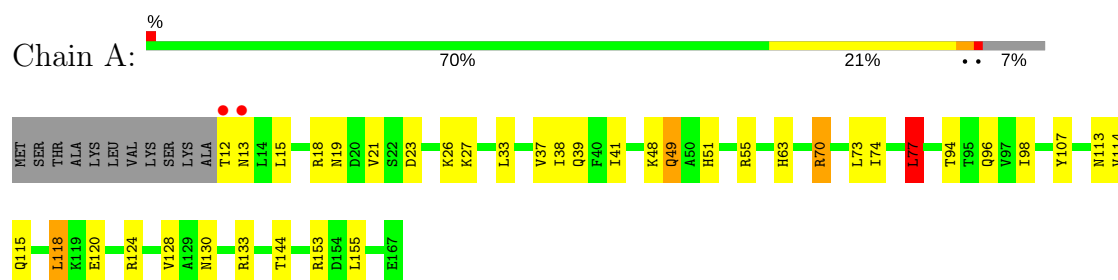
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	53	Total 53	O 53	0	0
3	B	36	Total 36	O 36	0	0
3	C	41	Total 41	O 41	0	0
3	D	45	Total 45	O 45	0	0
3	E	50	Total 50	O 50	0	0
3	F	41	Total 41	O 41	0	0
3	G	22	Total 22	O 22	0	0
3	H	19	Total 19	O 19	0	0
3	I	26	Total 26	O 26	0	0
3	J	17	Total 17	O 17	0	0
3	K	27	Total 27	O 27	0	0
3	L	9	Total 9	O 9	0	0

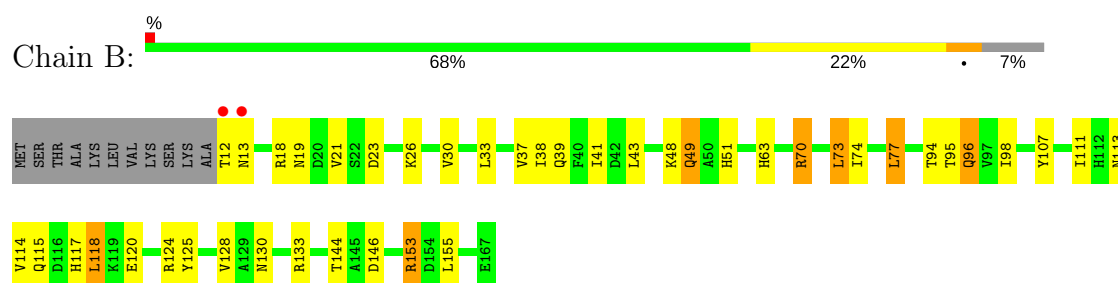
3 Residue-property plots [i](#)

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.

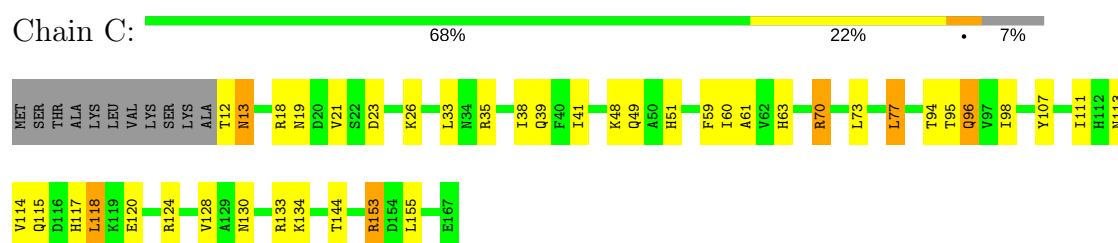
• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN



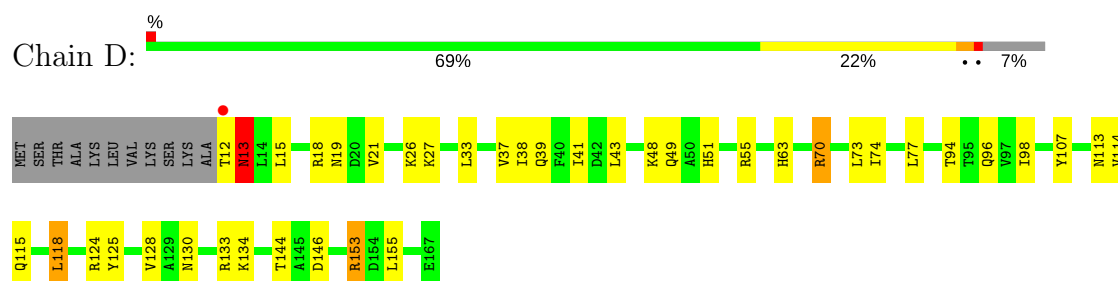
• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN



• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

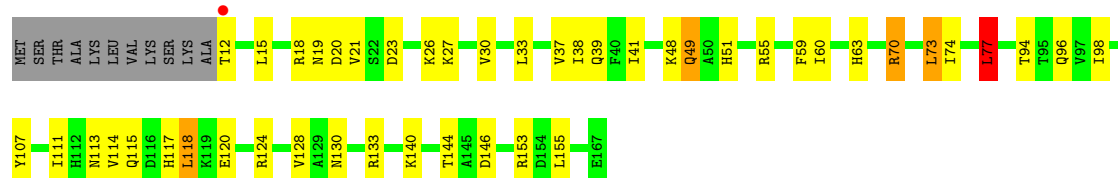


• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN



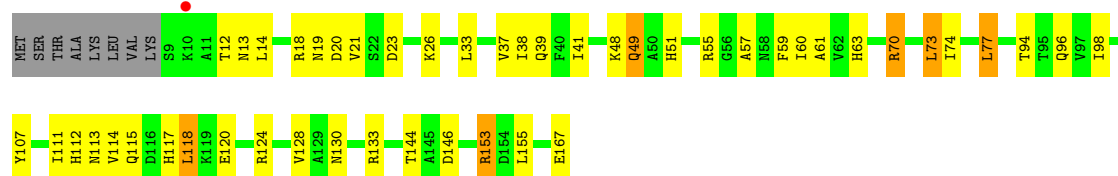
- Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

Chain E: 



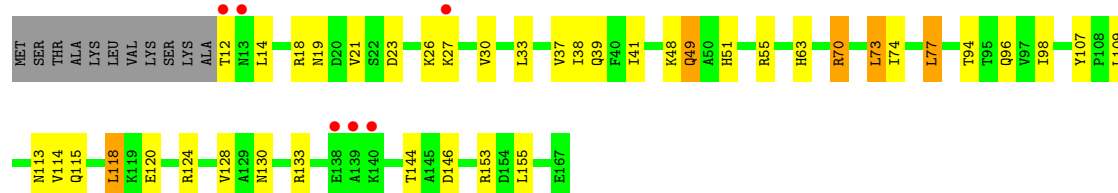
- Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

Chain F: 



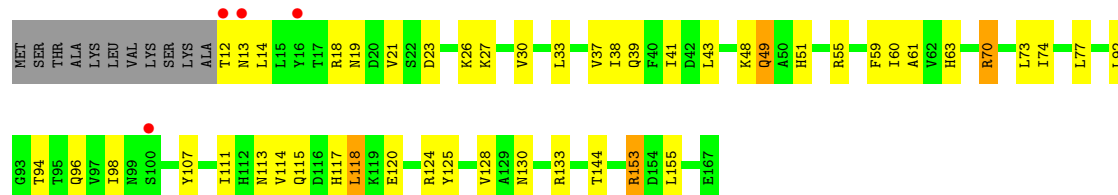
- Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

Chain G: 



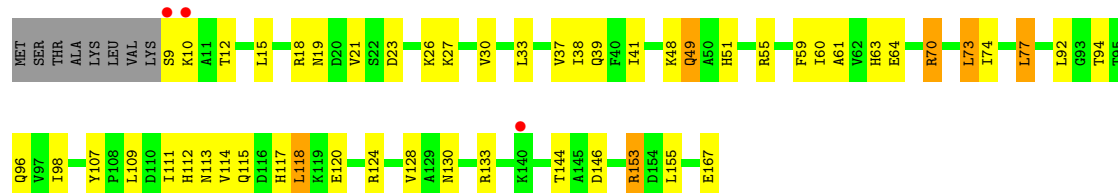
- Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

Chain H: 

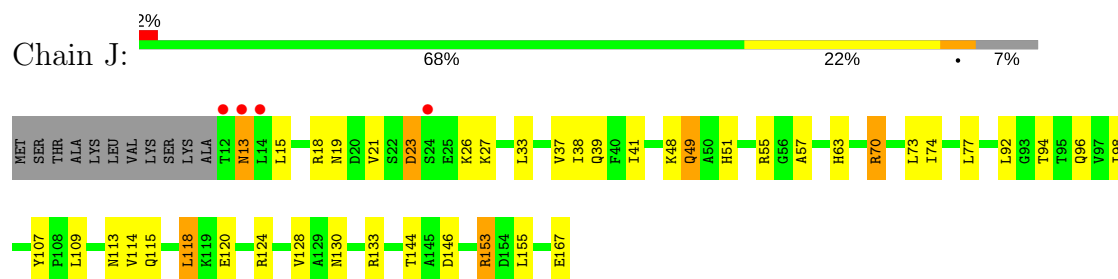


- Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN

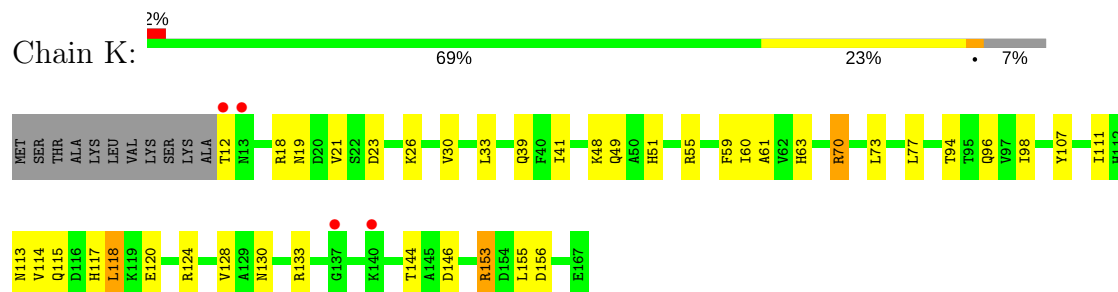
Chain I: 



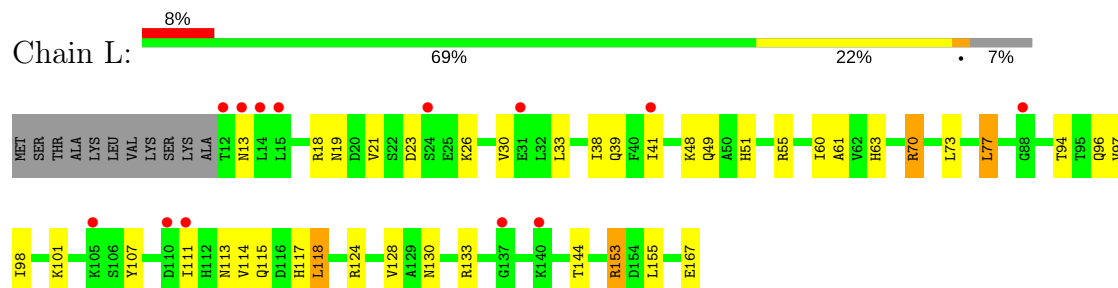
• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN



• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN



• Molecule 1: DNA PROTECTION DURING STARVATION PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	117.96Å 139.12Å 263.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.02 – 2.60 29.02 – 2.47	Depositor EDS
% Data completeness (in resolution range)	97.5 (29.02-2.60) 94.5 (29.02-2.47)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.01 (at 2.48Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.209 , 0.244 0.227 , 0.260	Depositor DCC
R_{free} test set	6588 reflections (10.13%)	DCC
Wilson B-factor (Å ²)	38.0	Xtriage
Anisotropy	0.132	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 42.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	15354	wwPDB-VP
Average B, all atoms (Å ²)	60.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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 ⓘ

5.1 Standard geometry ⓘ

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

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	0.37	0/1254	0.59	1/1698 (0.1%)
1	B	0.38	0/1254	0.60	1/1698 (0.1%)
1	C	0.38	0/1254	0.60	1/1698 (0.1%)
1	D	0.39	0/1254	0.59	0/1698
1	E	0.39	0/1254	0.60	1/1698 (0.1%)
1	F	0.38	0/1274	0.60	1/1724 (0.1%)
1	G	0.36	0/1254	0.57	1/1698 (0.1%)
1	H	0.34	0/1254	0.55	0/1698
1	I	0.34	0/1274	0.57	1/1724 (0.1%)
1	J	0.35	0/1254	0.57	0/1698
1	K	0.36	0/1254	0.58	0/1698
1	L	0.31	0/1254	0.55	1/1698 (0.1%)
All	All	0.36	0/15088	0.58	8/20428 (0.0%)

There are no bond length outliers.

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	E	77	LEU	CA-CB-CG	5.97	129.03	115.30
1	A	77	LEU	CA-CB-CG	5.54	128.04	115.30
1	B	77	LEU	CA-CB-CG	5.41	127.75	115.30
1	C	77	LEU	CA-CB-CG	5.36	127.62	115.30
1	G	77	LEU	CA-CB-CG	5.34	127.59	115.30
1	F	77	LEU	CA-CB-CG	5.23	127.32	115.30
1	L	77	LEU	CA-CB-CG	5.04	126.88	115.30
1	I	77	LEU	CA-CB-CG	5.01	126.82	115.30

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	1236	0	1232	35	0
1	B	1236	0	1232	39	0
1	C	1236	0	1232	31	0
1	D	1236	0	1232	36	0
1	E	1236	0	1232	43	0
1	F	1256	0	1255	46	0
1	G	1236	0	1232	37	0
1	H	1236	0	1232	43	0
1	I	1256	0	1255	46	0
1	J	1236	0	1232	40	0
1	K	1236	0	1232	28	0
1	L	1236	0	1232	28	0
2	A	8	0	12	0	0
2	C	16	0	24	2	0
2	E	8	0	12	0	0
2	F	16	0	24	4	0
2	G	8	0	12	3	0
2	I	8	0	12	0	0
2	J	8	0	12	1	0
2	K	16	0	24	1	0
2	L	8	0	12	2	0
3	A	53	0	0	1	0
3	B	36	0	0	0	0
3	C	41	0	0	3	0
3	D	45	0	0	1	0
3	E	50	0	0	1	0
3	F	41	0	0	1	0
3	G	22	0	0	1	0
3	H	19	0	0	0	0
3	I	26	0	0	0	0
3	J	17	0	0	0	0
3	K	27	0	0	1	0
3	L	9	0	0	0	0
All	All	15354	0	14974	416	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 14.

All (416) 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:J:41:ILE:HG12	1:J:77:LEU:HD11	1.49	0.94
1:H:41:ILE:HG12	1:H:77:LEU:HD11	1.55	0.87
1:L:41:ILE:HG12	1:L:77:LEU:HD11	1.57	0.85
1:B:41:ILE:HG12	1:B:77:LEU:HD11	1.59	0.85
1:D:41:ILE:HG12	1:D:77:LEU:HD11	1.59	0.84
1:G:41:ILE:HG12	1:G:77:LEU:HD11	1.60	0.83
1:A:41:ILE:HG12	1:A:77:LEU:HD11	1.60	0.82
1:F:41:ILE:HG12	1:F:77:LEU:HD11	1.60	0.82
1:I:41:ILE:HG12	1:I:77:LEU:HD11	1.60	0.81
1:C:12:THR:O	1:C:13:ASN:HB2	1.80	0.80
1:K:41:ILE:HG12	1:K:77:LEU:HD11	1.65	0.79
1:E:41:ILE:HG12	1:E:77:LEU:HD11	1.64	0.78
2:C:603:TRS:H22	1:E:146:ASP:OD1	1.85	0.75
1:G:94:THR:HG23	1:H:49:GLN:HG3	1.70	0.73
1:E:94:THR:HG23	1:F:49:GLN:HG3	1.70	0.72
1:I:94:THR:HG23	1:J:49:GLN:HG3	1.71	0.72
1:C:41:ILE:HG12	1:C:77:LEU:HD11	1.71	0.72
1:I:49:GLN:HG3	1:J:94:THR:HG23	1.72	0.72
1:B:39:GLN:HG2	1:B:128:VAL:HG22	1.74	0.70
1:E:49:GLN:HG3	1:F:94:THR:HG23	1.74	0.70
1:D:39:GLN:HG2	1:D:128:VAL:HG22	1.75	0.69
1:A:94:THR:HG23	1:B:49:GLN:HG3	1.74	0.68
1:J:113:ASN:ND2	1:J:115:GLN:HB2	2.10	0.67
1:A:21:VAL:O	1:A:26:LYS:HE3	1.94	0.67
1:G:14:LEU:HG	1:G:27:LYS:HE3	1.76	0.66
1:K:113:ASN:ND2	1:K:115:GLN:HB2	2.11	0.65
1:D:18:ARG:HG3	1:D:18:ARG:HH21	1.61	0.65
1:C:33:LEU:HD21	1:C:144:THR:HG23	1.77	0.65
1:E:94:THR:CG2	1:F:49:GLN:HG3	2.27	0.65
1:A:49:GLN:HG3	1:B:94:THR:HG23	1.76	0.65
1:H:33:LEU:HD21	1:H:144:THR:HG23	1.78	0.65
1:H:21:VAL:O	1:H:26:LYS:HE3	1.97	0.65
1:B:113:ASN:ND2	1:B:115:GLN:HB2	2.11	0.65
1:F:39:GLN:HG2	1:F:128:VAL:HG22	1.80	0.64
1:J:39:GLN:HG2	1:J:128:VAL:HG22	1.79	0.64
1:G:21:VAL:O	1:G:26:LYS:HE3	1.97	0.64
1:L:21:VAL:O	1:L:26:LYS:HE3	1.98	0.64
1:D:21:VAL:O	1:D:26:LYS:HE3	1.98	0.63
1:K:39:GLN:HG2	1:K:128:VAL:HG22	1.80	0.63
1:J:21:VAL:O	1:J:26:LYS:HE3	1.99	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:13:ASN:HB3	1:A:27:LYS:NZ	2.14	0.63
1:G:33:LEU:HD21	1:G:144:THR:HG23	1.80	0.63
1:H:113:ASN:ND2	1:H:115:GLN:HB2	2.14	0.63
1:G:49:GLN:HG3	1:H:94:THR:HG23	1.79	0.62
1:L:33:LEU:HD21	1:L:144:THR:HG23	1.81	0.62
1:I:113:ASN:ND2	1:I:115:GLN:HB2	2.14	0.62
1:L:113:ASN:ND2	1:L:115:GLN:HB2	2.14	0.62
1:E:12:THR:HG23	1:E:27:LYS:HG2	1.79	0.62
1:I:21:VAL:O	1:I:26:LYS:HE3	1.98	0.62
1:K:21:VAL:O	1:K:26:LYS:HE3	1.99	0.62
1:B:33:LEU:HD21	1:B:144:THR:HG23	1.81	0.62
1:E:49:GLN:HG3	1:F:94:THR:CG2	2.28	0.62
1:I:18:ARG:HH21	1:I:18:ARG:HG3	1.64	0.62
1:H:13:ASN:HB3	1:H:27:LYS:HE2	1.81	0.62
1:L:97:VAL:HG13	1:L:101:LYS:HE2	1.82	0.62
1:E:33:LEU:HD21	1:E:144:THR:HG23	1.82	0.61
1:C:39:GLN:HG2	1:C:128:VAL:HG22	1.81	0.61
1:A:39:GLN:HG2	1:A:128:VAL:HG22	1.83	0.61
1:F:33:LEU:HD21	1:F:144:THR:HG23	1.81	0.61
1:A:113:ASN:ND2	1:A:115:GLN:HB2	2.15	0.60
1:G:94:THR:CG2	1:H:49:GLN:HG3	2.31	0.60
1:L:39:GLN:HG2	1:L:128:VAL:HG22	1.83	0.60
1:H:39:GLN:HG2	1:H:128:VAL:HG22	1.82	0.60
1:E:48:LYS:HE3	1:E:70:ARG:NH1	2.16	0.60
1:K:48:LYS:HE3	1:K:70:ARG:NH1	2.15	0.60
1:A:94:THR:CG2	1:B:49:GLN:HG3	2.30	0.60
1:D:134:LYS:NZ	1:F:20:ASP:OD1	2.25	0.60
1:G:113:ASN:ND2	1:G:115:GLN:HB2	2.17	0.59
1:I:39:GLN:HG2	1:I:128:VAL:HG22	1.84	0.59
1:K:33:LEU:HD21	1:K:144:THR:HG23	1.84	0.59
1:C:113:ASN:ND2	1:C:115:GLN:HB2	2.18	0.59
1:E:39:GLN:HG2	1:E:128:VAL:HG22	1.83	0.59
1:D:33:LEU:HD21	1:D:144:THR:HG23	1.84	0.59
1:D:113:ASN:ND2	1:D:115:GLN:HB2	2.18	0.58
1:I:49:GLN:HG3	1:J:94:THR:CG2	2.32	0.58
1:F:113:ASN:ND2	1:F:115:GLN:HB2	2.18	0.58
1:I:94:THR:CG2	1:J:49:GLN:HG3	2.33	0.58
1:F:48:LYS:HE3	1:F:70:ARG:NH1	2.19	0.58
1:B:21:VAL:O	1:B:26:LYS:HE3	2.03	0.58
1:C:21:VAL:O	1:C:26:LYS:HE3	2.04	0.58
1:H:107:TYR:HD1	1:H:124:ARG:NH2	2.02	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:107:TYR:HD1	1:E:124:ARG:NH2	2.01	0.58
2:G:607:TRS:H12	1:K:146:ASP:OD1	2.03	0.57
1:I:48:LYS:HE3	1:I:70:ARG:NH1	2.19	0.57
1:F:21:VAL:O	1:F:26:LYS:HE3	2.03	0.57
1:E:113:ASN:ND2	1:E:115:GLN:HB2	2.19	0.57
1:F:18:ARG:HH21	1:F:18:ARG:HG3	1.69	0.57
1:E:18:ARG:HH21	1:E:18:ARG:HG3	1.69	0.57
1:J:33:LEU:HD21	1:J:144:THR:HG23	1.85	0.57
1:J:48:LYS:HE3	1:J:70:ARG:NH1	2.19	0.57
1:A:18:ARG:HG3	1:A:18:ARG:HH21	1.69	0.56
1:F:107:TYR:HD1	1:F:124:ARG:NH2	2.02	0.56
1:B:48:LYS:HE3	1:B:70:ARG:NH1	2.20	0.56
1:J:113:ASN:HD21	1:J:115:GLN:HB2	1.68	0.56
1:K:107:TYR:HD1	1:K:124:ARG:NH2	2.02	0.56
1:C:18:ARG:HH21	1:C:18:ARG:HG3	1.70	0.56
1:G:48:LYS:HE3	1:G:70:ARG:NH1	2.21	0.56
1:A:48:LYS:HE3	1:A:70:ARG:NH1	2.21	0.56
1:H:18:ARG:HH21	1:H:18:ARG:HG3	1.69	0.56
1:G:39:GLN:HG2	1:G:128:VAL:HG22	1.86	0.56
1:A:33:LEU:HD21	1:A:144:THR:HG23	1.88	0.56
1:L:18:ARG:HG3	1:L:18:ARG:HH21	1.71	0.56
1:L:48:LYS:HE3	1:L:70:ARG:NH1	2.21	0.55
1:A:13:ASN:HB3	1:A:27:LYS:HZ1	1.72	0.55
1:E:21:VAL:O	1:E:26:LYS:HE3	2.05	0.55
1:A:49:GLN:HG3	1:B:94:THR:CG2	2.36	0.55
1:H:48:LYS:HE3	1:H:70:ARG:NH1	2.21	0.55
1:J:18:ARG:HH21	1:J:18:ARG:HG3	1.72	0.55
1:F:70:ARG:NH2	3:F:618:HOH:O	2.39	0.55
1:K:130:ASN:OD1	1:K:133:ARG:NH2	2.39	0.55
1:F:12:THR:HG23	1:F:12:THR:O	2.05	0.55
1:C:107:TYR:HD1	1:C:124:ARG:NH2	2.06	0.54
1:G:18:ARG:HH21	1:G:18:ARG:HG3	1.72	0.54
1:I:33:LEU:HD21	1:I:144:THR:HG23	1.88	0.54
1:L:113:ASN:HD21	1:L:115:GLN:HB2	1.73	0.54
1:G:49:GLN:HG3	1:H:94:THR:CG2	2.38	0.54
1:B:113:ASN:HD21	1:B:115:GLN:HB2	1.74	0.53
1:I:107:TYR:HD1	1:I:124:ARG:NH2	2.05	0.53
1:E:130:ASN:OD1	1:E:133:ARG:NH2	2.39	0.53
1:C:48:LYS:HE3	1:C:70:ARG:NH1	2.24	0.53
1:I:74:ILE:HG23	1:J:70:ARG:NH2	2.24	0.53
1:H:113:ASN:HD21	1:H:115:GLN:HB2	1.73	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:107:TYR:HD1	1:D:124:ARG:NH2	2.06	0.53
1:G:70:ARG:NH2	1:H:74:ILE:HG23	2.24	0.53
1:B:107:TYR:HD1	1:B:124:ARG:NH2	2.07	0.52
1:G:146:ASP:OD1	2:G:607:TRS:H22	2.09	0.52
1:L:97:VAL:HG13	1:L:101:LYS:CE	2.39	0.52
1:C:51:HIS:CE1	1:C:63:HIS:CE1	2.98	0.52
1:D:130:ASN:OD1	1:D:133:ARG:NH2	2.39	0.52
1:E:15:LEU:N	1:E:15:LEU:HD12	2.24	0.52
1:J:130:ASN:OD1	1:J:133:ARG:NH2	2.42	0.52
1:G:107:TYR:HD1	1:G:124:ARG:NH2	2.07	0.52
1:I:51:HIS:CE1	1:I:63:HIS:CE1	2.98	0.52
1:L:107:TYR:HD1	1:L:124:ARG:NH2	2.08	0.51
1:B:18:ARG:HH21	1:B:18:ARG:HG3	1.75	0.51
1:B:39:GLN:CG	1:B:128:VAL:HG22	2.39	0.51
1:K:113:ASN:HD21	1:K:115:GLN:HB2	1.71	0.51
1:A:19:ASN:OD1	1:A:21:VAL:HG23	2.11	0.51
1:A:70:ARG:NH2	1:B:74:ILE:HG23	2.25	0.51
1:I:109:LEU:HD13	1:J:92:LEU:HD22	1.92	0.51
1:I:113:ASN:HD21	1:I:115:GLN:HB2	1.75	0.51
1:D:18:ARG:HG3	1:D:18:ARG:NH2	2.26	0.51
2:G:607:TRS:H32	1:I:146:ASP:OD1	2.11	0.51
1:D:153:ARG:HG3	3:D:342:HOH:O	2.09	0.51
1:F:55:ARG:O	1:F:114:VAL:HG23	2.11	0.51
1:G:19:ASN:OD1	1:G:21:VAL:HG23	2.11	0.50
1:J:57:ALA:HA	3:K:618:HOH:O	2.11	0.50
1:D:39:GLN:CG	1:D:128:VAL:HG22	2.41	0.50
1:E:70:ARG:NH2	1:F:74:ILE:HG23	2.27	0.50
1:B:12:THR:O	1:B:13:ASN:HB2	2.12	0.50
1:C:130:ASN:OD1	1:C:133:ARG:NH2	2.43	0.50
1:D:12:THR:N	1:D:27:LYS:HE2	2.27	0.50
1:G:130:ASN:OD1	1:G:133:ARG:NH2	2.44	0.50
1:K:18:ARG:HH21	1:K:18:ARG:HG3	1.77	0.49
1:A:74:ILE:HG23	1:B:70:ARG:NH2	2.27	0.49
1:A:113:ASN:HD21	1:A:115:GLN:HB2	1.78	0.49
1:C:153:ARG:NH1	1:C:153:ARG:HG2	2.27	0.49
1:D:51:HIS:CE1	1:D:63:HIS:CE1	3.00	0.49
1:A:37:VAL:O	1:A:41:ILE:HG13	2.11	0.49
1:B:114:VAL:HG12	1:B:118:LEU:HD22	1.94	0.49
1:I:153:ARG:NH1	1:I:153:ARG:HG2	2.26	0.49
1:F:130:ASN:OD1	1:F:133:ARG:NH2	2.43	0.49
1:C:153:ARG:HG2	1:C:153:ARG:HH11	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:130:ASN:OD1	1:A:133:ARG:NH2	2.44	0.49
1:C:39:GLN:CG	1:C:128:VAL:HG22	2.43	0.49
1:F:113:ASN:HD21	1:F:115:GLN:HB2	1.76	0.49
1:I:70:ARG:NH2	1:J:74:ILE:HG23	2.26	0.49
1:K:51:HIS:CE1	1:K:63:HIS:CE1	3.00	0.49
1:K:120:GLU:O	1:K:124:ARG:HG2	2.12	0.49
1:B:19:ASN:OD1	1:B:21:VAL:HG23	2.13	0.48
1:F:12:THR:C	1:F:14:LEU:H	2.16	0.48
1:L:130:ASN:OD1	1:L:133:ARG:NH2	2.46	0.48
1:I:153:ARG:HH11	1:I:153:ARG:HG2	1.77	0.48
1:C:113:ASN:HD21	1:C:115:GLN:HB2	1.76	0.48
1:D:12:THR:O	1:D:13:ASN:HB2	2.14	0.48
1:G:14:LEU:HG	1:G:27:LYS:CE	2.44	0.48
1:A:39:GLN:CG	1:A:128:VAL:HG22	2.43	0.48
1:D:48:LYS:HE3	1:D:70:ARG:NH1	2.28	0.48
1:J:114:VAL:HG12	1:J:118:LEU:HD22	1.96	0.48
1:H:51:HIS:CE1	1:H:63:HIS:CE1	3.02	0.48
1:G:109:LEU:HD13	1:H:92:LEU:HD22	1.96	0.48
1:H:94:THR:O	1:H:98:ILE:HG12	2.13	0.48
1:J:19:ASN:OD1	1:J:21:VAL:HG23	2.14	0.48
1:J:51:HIS:CE1	1:J:63:HIS:CE1	3.02	0.48
1:D:146:ASP:OD2	2:F:606:TRS:H12	2.13	0.48
1:G:37:VAL:O	1:G:41:ILE:HG13	2.14	0.48
1:B:51:HIS:CE1	1:B:63:HIS:CE1	3.01	0.47
1:H:39:GLN:CG	1:H:128:VAL:HG22	2.44	0.47
1:I:55:ARG:O	1:I:114:VAL:HG23	2.14	0.47
1:E:114:VAL:HG12	1:E:118:LEU:HD22	1.96	0.47
1:F:19:ASN:OD1	1:F:21:VAL:HG23	2.15	0.47
1:G:38:ILE:HG23	1:G:98:ILE:HA	1.96	0.47
1:I:18:ARG:HG3	1:I:18:ARG:NH2	2.29	0.47
1:A:107:TYR:HD1	1:A:124:ARG:NH2	2.12	0.47
1:E:111:ILE:HG13	1:E:117:HIS:CE1	2.49	0.47
1:I:37:VAL:O	1:I:41:ILE:HG13	2.15	0.47
1:K:94:THR:O	1:K:98:ILE:HG12	2.15	0.47
1:F:18:ARG:HG3	1:F:18:ARG:NH2	2.29	0.47
1:G:113:ASN:HD21	1:G:115:GLN:HB2	1.78	0.47
1:L:94:THR:O	1:L:98:ILE:HG12	2.15	0.47
1:C:19:ASN:OD1	1:C:21:VAL:HG23	2.15	0.47
1:H:130:ASN:OD1	1:H:133:ARG:NH2	2.45	0.47
1:K:153:ARG:NH1	1:K:153:ARG:HG2	2.29	0.47
1:J:167:GLU:OE1	2:K:611:TRS:H21	2.15	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:55:ARG:O	1:D:114:VAL:HG23	2.14	0.47
1:I:120:GLU:O	1:I:124:ARG:HG2	2.14	0.47
1:I:9:SER:OG	1:I:10:LYS:N	2.47	0.47
1:C:120:GLU:O	1:C:124:ARG:HG2	2.15	0.47
1:D:153:ARG:HG2	1:D:153:ARG:NH1	2.30	0.47
1:I:74:ILE:HG23	1:J:70:ARG:HH22	1.80	0.47
1:I:38:ILE:HG23	1:I:98:ILE:HA	1.97	0.47
1:A:51:HIS:CE1	1:A:63:HIS:CE1	3.03	0.47
1:I:94:THR:O	1:I:98:ILE:HG12	2.15	0.47
1:K:39:GLN:CG	1:K:128:VAL:HG22	2.43	0.46
1:F:51:HIS:CE1	1:F:63:HIS:CE1	3.03	0.46
1:G:94:THR:O	1:G:98:ILE:HG12	2.16	0.46
1:H:37:VAL:O	1:H:41:ILE:HG13	2.15	0.46
1:J:107:TYR:HD1	1:J:124:ARG:NH2	2.13	0.46
1:G:55:ARG:O	1:G:114:VAL:HG23	2.15	0.46
1:L:19:ASN:OD1	1:L:21:VAL:HG23	2.15	0.46
1:H:26:LYS:O	1:H:30:VAL:HG23	2.14	0.46
1:D:38:ILE:HG23	1:D:98:ILE:HA	1.96	0.46
1:F:37:VAL:O	1:F:41:ILE:HG13	2.16	0.46
1:F:60:ILE:HG23	1:F:61:ALA:N	2.31	0.46
1:F:12:THR:CG2	1:F:12:THR:O	2.64	0.46
1:I:60:ILE:O	1:I:64:GLU:HG3	2.16	0.46
1:K:107:TYR:CD1	1:K:124:ARG:NH2	2.83	0.46
1:E:51:HIS:CE1	1:E:63:HIS:CE1	3.03	0.46
1:L:167:GLU:OE1	2:L:608:TRS:H31	2.16	0.46
1:B:130:ASN:OD1	1:B:133:ARG:NH2	2.43	0.46
1:I:112:HIS:CE1	1:J:15:LEU:HD21	2.51	0.46
1:I:19:ASN:OD1	1:I:21:VAL:HG23	2.16	0.45
1:I:48:LYS:HA	1:I:51:HIS:HB3	1.98	0.45
1:J:55:ARG:O	1:J:114:VAL:HG23	2.16	0.45
1:L:39:GLN:CG	1:L:128:VAL:HG22	2.45	0.45
1:B:153:ARG:NH1	1:B:153:ARG:HG2	2.30	0.45
1:H:14:LEU:HD11	1:H:27:LYS:HG3	1.98	0.45
1:J:146:ASP:OD1	2:J:612:TRS:H32	2.17	0.45
1:D:12:THR:O	1:D:13:ASN:CB	2.65	0.45
1:H:153:ARG:HG2	1:H:153:ARG:NH1	2.31	0.45
1:K:118:LEU:HA	1:K:118:LEU:HD12	1.85	0.45
1:E:19:ASN:OD1	1:E:21:VAL:HG23	2.16	0.45
1:G:114:VAL:HG12	1:G:118:LEU:HD22	1.98	0.45
1:E:48:LYS:HA	1:E:51:HIS:HB3	1.98	0.45
1:F:167:GLU:OE1	2:F:605:TRS:H21	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:48:LYS:HA	1:G:51:HIS:HB3	1.99	0.45
1:I:114:VAL:HG12	1:I:118:LEU:HD22	1.97	0.45
1:J:48:LYS:HA	1:J:51:HIS:HB3	1.99	0.45
1:J:94:THR:O	1:J:98:ILE:HG12	2.16	0.45
1:A:114:VAL:HG12	1:A:118:LEU:HD22	1.98	0.45
1:J:13:ASN:HB3	1:J:27:LYS:NZ	2.32	0.45
1:C:94:THR:O	1:C:98:ILE:HG12	2.17	0.45
1:K:114:VAL:HG12	1:K:118:LEU:HD22	1.98	0.45
3:C:624:HOH:O	1:F:57:ALA:HA	2.16	0.45
1:I:60:ILE:HG23	1:I:61:ALA:N	2.32	0.45
1:A:94:THR:O	1:A:98:ILE:HG12	2.16	0.45
1:D:114:VAL:HG12	1:D:118:LEU:HD22	1.99	0.45
1:F:94:THR:O	1:F:98:ILE:HG12	2.17	0.45
1:J:38:ILE:HG23	1:J:98:ILE:HA	1.98	0.45
1:C:18:ARG:HG3	1:C:18:ARG:NH2	2.32	0.44
1:D:43:LEU:HD11	1:D:125:TYR:CD1	2.52	0.44
1:F:48:LYS:HA	1:F:51:HIS:HB3	2.00	0.44
1:L:51:HIS:CE1	1:L:63:HIS:CE1	3.06	0.44
1:G:120:GLU:O	1:G:124:ARG:HG2	2.16	0.44
1:G:39:GLN:CG	1:G:128:VAL:HG22	2.46	0.44
1:G:51:HIS:CE1	1:G:63:HIS:CE1	3.05	0.44
1:G:55:ARG:NE	3:G:614:HOH:O	2.48	0.44
1:K:153:ARG:HH11	1:K:153:ARG:HG2	1.82	0.44
1:B:153:ARG:HH11	1:B:153:ARG:HG2	1.82	0.44
1:D:113:ASN:HD21	1:D:115:GLN:HB2	1.81	0.44
1:F:107:TYR:CD1	1:F:124:ARG:NH2	2.85	0.44
1:G:26:LYS:O	1:G:30:VAL:HG23	2.18	0.44
1:I:111:ILE:HG13	1:I:117:HIS:CE1	2.52	0.44
1:B:38:ILE:HG23	1:B:98:ILE:HA	1.99	0.44
1:F:38:ILE:HG23	1:F:98:ILE:HA	2.00	0.44
1:G:74:ILE:HG23	1:H:70:ARG:NH2	2.32	0.44
1:H:18:ARG:NH2	1:H:18:ARG:HG3	2.31	0.44
1:C:153:ARG:HG3	3:C:633:HOH:O	2.17	0.44
1:E:39:GLN:CG	1:E:128:VAL:HG22	2.47	0.44
1:B:95:THR:HG23	1:B:96:GLN:N	2.33	0.44
1:C:35:ARG:HD3	3:C:607:HOH:O	2.17	0.44
1:G:113:ASN:ND2	1:G:115:GLN:H	2.15	0.44
1:H:19:ASN:OD1	1:H:21:VAL:HG23	2.17	0.44
1:B:111:ILE:HG13	1:B:117:HIS:CE1	2.53	0.44
1:B:120:GLU:O	1:B:124:ARG:HG2	2.18	0.44
1:K:60:ILE:HG23	1:K:61:ALA:N	2.32	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:118:LEU:HA	1:A:118:LEU:HD12	1.90	0.44
1:C:48:LYS:HA	1:C:51:HIS:HB3	1.99	0.44
1:J:23:ASP:O	1:J:27:LYS:HG2	2.17	0.44
1:L:48:LYS:HA	1:L:51:HIS:HB3	2.00	0.44
1:D:19:ASN:OD1	1:D:21:VAL:HG23	2.18	0.43
1:D:37:VAL:O	1:D:41:ILE:HG13	2.18	0.43
1:I:92:LEU:HD22	1:J:109:LEU:HD13	2.00	0.43
1:L:118:LEU:HA	1:L:118:LEU:HD12	1.84	0.43
1:H:107:TYR:CD1	1:H:124:ARG:NH2	2.85	0.43
1:I:130:ASN:OD1	1:I:133:ARG:NH2	2.48	0.43
1:B:37:VAL:O	1:B:41:ILE:HG13	2.19	0.43
1:D:13:ASN:HD22	1:D:13:ASN:HA	1.57	0.43
1:D:153:ARG:HG2	1:D:153:ARG:HH11	1.81	0.43
1:C:134:LYS:NZ	1:E:20:ASP:OD1	2.35	0.43
1:F:39:GLN:CG	1:F:128:VAL:HG22	2.47	0.43
1:J:39:GLN:CG	1:J:128:VAL:HG22	2.46	0.43
1:L:114:VAL:HG12	1:L:118:LEU:HD22	1.99	0.43
1:D:94:THR:O	1:D:98:ILE:HG12	2.19	0.43
1:A:18:ARG:HG3	1:A:18:ARG:NH2	2.33	0.43
1:C:118:LEU:HD12	1:C:118:LEU:HA	1.88	0.43
1:C:153:ARG:CG	1:C:153:ARG:HH11	2.32	0.43
1:H:55:ARG:O	1:H:114:VAL:HG23	2.18	0.43
1:K:55:ARG:O	1:K:114:VAL:HG23	2.18	0.43
1:A:15:LEU:HD12	1:A:15:LEU:N	2.33	0.43
1:H:120:GLU:O	1:H:124:ARG:HG2	2.19	0.43
1:K:26:LYS:O	1:K:30:VAL:HG23	2.18	0.43
1:E:113:ASN:HD21	1:E:115:GLN:HB2	1.80	0.43
1:E:12:THR:O	1:E:12:THR:HG23	2.19	0.43
1:I:26:LYS:O	1:I:30:VAL:HG23	2.19	0.43
1:K:111:ILE:HG13	1:K:117:HIS:CE1	2.54	0.43
1:K:19:ASN:OD1	1:K:21:VAL:HG23	2.18	0.43
1:E:38:ILE:HG23	1:E:98:ILE:HA	2.01	0.43
1:F:113:ASN:ND2	1:F:115:GLN:H	2.16	0.43
1:G:18:ARG:NH2	1:G:18:ARG:HG3	2.34	0.43
1:H:48:LYS:HA	1:H:51:HIS:HB3	2.01	0.43
1:A:55:ARG:O	1:A:114:VAL:HG23	2.19	0.42
1:D:118:LEU:HA	1:D:118:LEU:HD12	1.92	0.42
1:E:55:ARG:O	1:E:114:VAL:HG23	2.18	0.42
1:F:70:ARG:HD3	1:F:74:ILE:HD12	2.01	0.42
1:C:59:PHE:CD1	1:C:60:ILE:N	2.87	0.42
1:H:153:ARG:HH11	1:H:153:ARG:HG2	1.84	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:12:THR:O	1:I:27:LYS:HE3	2.19	0.42
1:L:18:ARG:NH2	1:L:18:ARG:HG3	2.33	0.42
1:B:146:ASP:OD1	2:C:603:TRS:O2	2.28	0.42
1:E:12:THR:CG2	1:E:27:LYS:HG2	2.45	0.42
1:I:15:LEU:HD12	1:I:15:LEU:N	2.33	0.42
1:K:133:ARG:NH1	1:K:156:ASP:OD2	2.52	0.42
1:A:12:THR:O	1:A:13:ASN:HB3	2.20	0.42
1:E:59:PHE:CD1	1:E:60:ILE:N	2.87	0.42
1:J:18:ARG:NH2	1:J:18:ARG:HG3	2.33	0.42
1:L:55:ARG:O	1:L:114:VAL:HG23	2.19	0.42
1:D:113:ASN:ND2	1:D:115:GLN:H	2.17	0.42
1:D:107:TYR:CD1	1:D:124:ARG:NH2	2.87	0.42
1:J:120:GLU:O	1:J:124:ARG:HG2	2.19	0.42
1:B:113:ASN:ND2	1:B:115:GLN:H	2.18	0.42
1:C:114:VAL:HG12	1:C:118:LEU:HD22	2.01	0.42
1:H:13:ASN:HB3	1:H:27:LYS:CE	2.50	0.42
1:G:70:ARG:HH22	1:H:74:ILE:HG23	1.84	0.42
1:I:59:PHE:CD1	1:I:60:ILE:N	2.88	0.42
1:B:48:LYS:HE3	1:B:70:ARG:CZ	2.50	0.42
1:D:15:LEU:N	1:D:15:LEU:HD12	2.34	0.42
1:E:74:ILE:HG23	1:F:70:ARG:NH2	2.34	0.42
1:G:73:LEU:HD12	1:G:73:LEU:HA	1.90	0.42
1:H:60:ILE:HG23	1:H:61:ALA:N	2.35	0.42
1:E:18:ARG:HG3	1:E:18:ARG:NH2	2.33	0.42
1:G:12:THR:HG23	1:G:12:THR:O	2.19	0.42
1:H:111:ILE:HG13	1:H:117:HIS:CE1	2.55	0.42
1:I:39:GLN:CG	1:I:128:VAL:HG22	2.48	0.42
1:K:77:LEU:C	1:K:77:LEU:HD23	2.40	0.42
1:L:38:ILE:HD13	1:L:101:LYS:HE3	2.02	0.42
1:B:70:ARG:HD3	1:B:74:ILE:HD12	2.02	0.42
1:C:111:ILE:HG13	1:C:117:HIS:CE1	2.55	0.42
1:E:94:THR:O	1:E:98:ILE:HG12	2.19	0.42
1:J:153:ARG:HG2	1:J:153:ARG:NH1	2.34	0.42
1:E:118:LEU:HD12	1:E:118:LEU:HA	1.91	0.41
1:E:107:TYR:CD1	1:E:124:ARG:NH2	2.85	0.41
1:H:43:LEU:HD11	1:H:125:TYR:CD1	2.55	0.41
1:L:111:ILE:HG13	1:L:117:HIS:CE1	2.55	0.41
1:L:60:ILE:HG23	1:L:61:ALA:N	2.35	0.41
1:E:120:GLU:O	1:E:124:ARG:HG2	2.20	0.41
1:E:70:ARG:HH22	1:F:74:ILE:HG23	1.86	0.41
1:A:74:ILE:HG23	1:B:70:ARG:HH22	1.86	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:38:ILE:HG23	1:C:98:ILE:HA	2.02	0.41
1:E:37:VAL:O	1:E:41:ILE:HG13	2.21	0.41
1:F:118:LEU:HD12	1:F:118:LEU:HA	1.89	0.41
1:I:70:ARG:HH22	1:J:74:ILE:HG23	1.84	0.41
1:D:70:ARG:HD3	1:D:74:ILE:HD12	2.01	0.41
1:L:26:LYS:O	1:L:30:VAL:HG23	2.20	0.41
1:L:38:ILE:HG23	1:L:98:ILE:HA	2.02	0.41
1:A:120:GLU:O	1:A:124:ARG:HG2	2.19	0.41
1:B:18:ARG:NH2	1:B:18:ARG:HG3	2.36	0.41
1:F:111:ILE:HG13	1:F:117:HIS:CE1	2.56	0.41
1:B:73:LEU:HD12	1:B:73:LEU:HA	1.92	0.41
1:H:77:LEU:C	1:H:77:LEU:HD23	2.40	0.41
1:H:38:ILE:HG23	1:H:98:ILE:HA	2.03	0.41
1:B:94:THR:O	1:B:98:ILE:HG12	2.21	0.41
1:E:26:LYS:O	1:E:30:VAL:HG23	2.20	0.41
1:I:153:ARG:HH11	1:I:153:ARG:CG	2.33	0.41
1:B:26:LYS:O	1:B:30:VAL:HG23	2.21	0.41
1:C:60:ILE:HG23	1:C:61:ALA:N	2.35	0.41
1:E:140:LYS:N	3:E:605:HOH:O	2.54	0.41
1:H:113:ASN:ND2	1:H:115:GLN:H	2.19	0.41
1:H:59:PHE:CD1	1:H:60:ILE:N	2.89	0.41
1:B:43:LEU:HD11	1:B:125:TYR:CD1	2.56	0.41
1:C:95:THR:HG23	1:C:96:GLN:N	2.34	0.41
1:F:73:LEU:HD12	1:F:73:LEU:HA	1.95	0.41
1:F:153:ARG:HG2	1:F:153:ARG:NH1	2.36	0.41
1:F:146:ASP:OD1	2:F:606:TRS:C3	2.69	0.41
1:I:73:LEU:HA	1:I:73:LEU:HD12	1.94	0.41
1:J:37:VAL:O	1:J:41:ILE:HG13	2.21	0.41
1:I:167:GLU:OE1	2:L:608:TRS:H21	2.21	0.41
1:A:13:ASN:CB	1:A:27:LYS:NZ	2.80	0.40
1:A:49:GLN:NE2	3:A:276:HOH:O	2.53	0.40
1:D:48:LYS:HA	1:D:51:HIS:HB3	2.02	0.40
1:F:120:GLU:O	1:F:124:ARG:HG2	2.21	0.40
1:F:167:GLU:CD	2:F:605:TRS:H21	2.40	0.40
1:J:153:ARG:HG2	1:J:153:ARG:HH11	1.86	0.40
1:A:48:LYS:HA	1:A:51:HIS:HB3	2.03	0.40
1:A:38:ILE:HG23	1:A:98:ILE:HA	2.04	0.40
1:E:15:LEU:HD21	1:F:112:HIS:CE1	2.57	0.40
1:E:48:LYS:HE3	1:E:70:ARG:CZ	2.52	0.40
1:H:114:VAL:HG12	1:H:118:LEU:HD22	2.03	0.40
1:H:12:THR:CG2	1:H:13:ASN:N	2.83	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:48:LYS:HA	1:B:51:HIS:HB3	2.02	0.40
1:L:153:ARG:NH1	1:L:153:ARG:HG2	2.36	0.40
1:D:77:LEU:C	1:D:77:LEU:HD23	2.42	0.40
1:E:73:LEU:HA	1:E:73:LEU:HD12	1.94	0.40
1:F:59:PHE:CD1	1:F:60:ILE:N	2.89	0.40
1:K:59:PHE:CD1	1:K:60:ILE:N	2.90	0.40
1:J:118:LEU:HA	1:J:118:LEU:HD12	1.90	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	154/167 (92%)	151 (98%)	3 (2%)	0	100	100
1	B	154/167 (92%)	151 (98%)	3 (2%)	0	100	100
1	C	154/167 (92%)	150 (97%)	3 (2%)	1 (1%)	28	53
1	D	154/167 (92%)	150 (97%)	3 (2%)	1 (1%)	28	53
1	E	154/167 (92%)	151 (98%)	3 (2%)	0	100	100
1	F	157/167 (94%)	153 (98%)	3 (2%)	1 (1%)	28	53
1	G	154/167 (92%)	151 (98%)	3 (2%)	0	100	100
1	H	154/167 (92%)	150 (97%)	4 (3%)	0	100	100
1	I	157/167 (94%)	152 (97%)	5 (3%)	0	100	100
1	J	154/167 (92%)	150 (97%)	3 (2%)	1 (1%)	28	53
1	K	154/167 (92%)	151 (98%)	3 (2%)	0	100	100
1	L	154/167 (92%)	151 (98%)	3 (2%)	0	100	100
All	All	1854/2004 (92%)	1811 (98%)	39 (2%)	4 (0%)	51	76

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	13	ASN
1	D	13	ASN
1	J	13	ASN
1	F	13	ASN

5.3.2 Protein sidechains ⓘ

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	134/143 (94%)	125 (93%)	9 (7%)	19	38
1	B	134/143 (94%)	126 (94%)	8 (6%)	22	44
1	C	134/143 (94%)	126 (94%)	8 (6%)	22	44
1	D	134/143 (94%)	126 (94%)	8 (6%)	22	44
1	E	134/143 (94%)	125 (93%)	9 (7%)	19	38
1	F	136/143 (95%)	128 (94%)	8 (6%)	23	45
1	G	134/143 (94%)	126 (94%)	8 (6%)	22	44
1	H	134/143 (94%)	126 (94%)	8 (6%)	22	44
1	I	136/143 (95%)	128 (94%)	8 (6%)	23	45
1	J	134/143 (94%)	126 (94%)	8 (6%)	22	44
1	K	134/143 (94%)	125 (93%)	9 (7%)	19	38
1	L	134/143 (94%)	125 (93%)	9 (7%)	19	38
All	All	1612/1716 (94%)	1512 (94%)	100 (6%)	21	42

All (100) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	ASP
1	A	49	GLN
1	A	70	ARG
1	A	73	LEU
1	A	77	LEU
1	A	96	GLN
1	A	118	LEU

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Mol	Chain	Res	Type
1	A	153	ARG
1	A	155	LEU
1	B	23	ASP
1	B	49	GLN
1	B	70	ARG
1	B	73	LEU
1	B	96	GLN
1	B	118	LEU
1	B	153	ARG
1	B	155	LEU
1	C	23	ASP
1	C	49	GLN
1	C	70	ARG
1	C	73	LEU
1	C	96	GLN
1	C	118	LEU
1	C	153	ARG
1	C	155	LEU
1	D	13	ASN
1	D	49	GLN
1	D	70	ARG
1	D	73	LEU
1	D	96	GLN
1	D	118	LEU
1	D	153	ARG
1	D	155	LEU
1	E	23	ASP
1	E	49	GLN
1	E	70	ARG
1	E	73	LEU
1	E	77	LEU
1	E	96	GLN
1	E	118	LEU
1	E	153	ARG
1	E	155	LEU
1	F	23	ASP
1	F	49	GLN
1	F	70	ARG
1	F	73	LEU
1	F	96	GLN
1	F	118	LEU
1	F	153	ARG

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Mol	Chain	Res	Type
1	F	155	LEU
1	G	23	ASP
1	G	49	GLN
1	G	70	ARG
1	G	73	LEU
1	G	96	GLN
1	G	118	LEU
1	G	153	ARG
1	G	155	LEU
1	H	23	ASP
1	H	49	GLN
1	H	70	ARG
1	H	73	LEU
1	H	96	GLN
1	H	118	LEU
1	H	153	ARG
1	H	155	LEU
1	I	23	ASP
1	I	49	GLN
1	I	70	ARG
1	I	73	LEU
1	I	96	GLN
1	I	118	LEU
1	I	153	ARG
1	I	155	LEU
1	J	23	ASP
1	J	49	GLN
1	J	70	ARG
1	J	73	LEU
1	J	96	GLN
1	J	118	LEU
1	J	153	ARG
1	J	155	LEU
1	K	12	THR
1	K	23	ASP
1	K	49	GLN
1	K	70	ARG
1	K	73	LEU
1	K	96	GLN
1	K	118	LEU
1	K	153	ARG
1	K	155	LEU

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Mol	Chain	Res	Type
1	L	13	ASN
1	L	23	ASP
1	L	49	GLN
1	L	70	ARG
1	L	73	LEU
1	L	96	GLN
1	L	118	LEU
1	L	153	ARG
1	L	155	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (58) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	36	GLN
1	A	39	GLN
1	A	96	GLN
1	A	113	ASN
1	B	13	ASN
1	B	36	GLN
1	B	39	GLN
1	B	96	GLN
1	B	113	ASN
1	C	36	GLN
1	C	49	GLN
1	C	96	GLN
1	C	113	ASN
1	D	13	ASN
1	D	36	GLN
1	D	39	GLN
1	D	49	GLN
1	D	96	GLN
1	D	113	ASN
1	E	36	GLN
1	E	39	GLN
1	E	96	GLN
1	E	113	ASN
1	F	36	GLN
1	F	39	GLN
1	F	96	GLN
1	F	99	ASN
1	F	112	HIS
1	F	113	ASN

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Mol	Chain	Res	Type
1	G	36	GLN
1	G	39	GLN
1	G	96	GLN
1	G	99	ASN
1	G	113	ASN
1	H	36	GLN
1	H	39	GLN
1	H	96	GLN
1	H	113	ASN
1	I	36	GLN
1	I	39	GLN
1	I	96	GLN
1	I	99	ASN
1	I	113	ASN
1	J	36	GLN
1	J	39	GLN
1	J	96	GLN
1	J	99	ASN
1	J	113	ASN
1	K	36	GLN
1	K	39	GLN
1	K	49	GLN
1	K	96	GLN
1	K	113	ASN
1	L	36	GLN
1	L	39	GLN
1	L	49	GLN
1	L	96	GLN
1	L	113	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 carbohydrates in this entry.

5.6 Ligand geometry

12 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	TRS	A	602	-	7,7,7	1.19	0	9,9,9	1.26	0
2	TRS	C	603	-	7,7,7	0.82	0	9,9,9	1.50	2 (22%)
2	TRS	C	604	-	7,7,7	1.18	1 (14%)	9,9,9	1.49	2 (22%)
2	TRS	E	601	-	7,7,7	0.62	0	9,9,9	1.45	1 (11%)
2	TRS	F	605	-	7,7,7	0.73	0	9,9,9	1.49	2 (22%)
2	TRS	F	606	-	7,7,7	0.81	0	9,9,9	1.42	1 (11%)
2	TRS	G	607	-	7,7,7	0.75	0	9,9,9	1.26	0
2	TRS	I	609	-	7,7,7	1.02	0	9,9,9	1.39	0
2	TRS	J	612	-	7,7,7	0.70	0	9,9,9	1.33	0
2	TRS	K	610	-	7,7,7	1.16	1 (14%)	9,9,9	1.40	1 (11%)
2	TRS	K	611	-	7,7,7	0.67	0	9,9,9	1.46	1 (11%)
2	TRS	L	608	-	7,7,7	0.68	0	9,9,9	1.49	1 (11%)

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	TRS	A	602	-	-	0/9/9/9	0/0/0/0
2	TRS	C	603	-	-	0/9/9/9	0/0/0/0
2	TRS	C	604	-	-	0/9/9/9	0/0/0/0
2	TRS	E	601	-	-	0/9/9/9	0/0/0/0
2	TRS	F	605	-	-	0/9/9/9	0/0/0/0
2	TRS	F	606	-	-	0/9/9/9	0/0/0/0
2	TRS	G	607	-	-	0/9/9/9	0/0/0/0
2	TRS	I	609	-	-	0/9/9/9	0/0/0/0
2	TRS	J	612	-	-	0/9/9/9	0/0/0/0
2	TRS	K	610	-	-	0/9/9/9	0/0/0/0
2	TRS	K	611	-	-	0/9/9/9	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	TRS	L	608	-	-	0/9/9/9	0/0/0/0

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	604	TRS	C2-C	2.07	1.57	1.52
2	K	610	TRS	C1-C	2.12	1.57	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	604	TRS	C2-C-C1	-2.17	104.91	111.06
2	F	605	TRS	C2-C-C1	-2.07	105.20	111.06
2	C	603	TRS	C2-C-N	2.04	112.07	107.73
2	K	611	TRS	C3-C-N	2.06	112.10	107.73
2	E	601	TRS	C3-C-N	2.12	112.24	107.73
2	L	608	TRS	C1-C-N	2.23	112.47	107.73
2	C	603	TRS	C1-C-N	2.26	112.53	107.73
2	F	605	TRS	C2-C-N	2.27	112.55	107.73
2	C	604	TRS	C2-C-N	2.42	112.88	107.73
2	K	610	TRS	C1-C-N	2.43	112.88	107.73
2	F	606	TRS	C1-C-N	2.72	113.50	107.73

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

7 monomers are involved in 13 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	603	TRS	2	0
2	F	605	TRS	2	0
2	F	606	TRS	2	0
2	G	607	TRS	3	0
2	J	612	TRS	1	0
2	K	611	TRS	1	0
2	L	608	TRS	2	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	156/167 (93%)	-0.50	2 (1%) 77 73	39, 50, 71, 99	0
1	B	156/167 (93%)	-0.35	2 (1%) 77 73	39, 52, 67, 88	0
1	C	156/167 (93%)	-0.37	0 100 100	38, 51, 71, 89	0
1	D	156/167 (93%)	-0.46	1 (0%) 89 88	36, 49, 68, 78	0
1	E	156/167 (93%)	-0.41	1 (0%) 89 88	35, 49, 66, 76	0
1	F	159/167 (95%)	-0.41	1 (0%) 89 88	40, 50, 71, 91	0
1	G	156/167 (93%)	-0.08	6 (3%) 41 33	47, 64, 87, 105	0
1	H	156/167 (93%)	0.09	4 (2%) 56 49	57, 72, 86, 94	0
1	I	159/167 (95%)	-0.17	3 (1%) 67 61	45, 65, 81, 87	0
1	J	156/167 (93%)	-0.07	4 (2%) 56 49	49, 64, 90, 108	0
1	K	156/167 (93%)	-0.14	4 (2%) 56 49	46, 61, 82, 111	0
1	L	156/167 (93%)	0.30	13 (8%) 12 8	58, 79, 94, 111	0
All	All	1878/2004 (93%)	-0.21	41 (2%) 62 56	35, 58, 85, 111	0

All (41) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	J	12	THR	5.4
1	G	12	THR	5.4
1	A	12	THR	5.2
1	H	12	THR	5.1
1	K	12	THR	5.0
1	L	12	THR	5.0
1	L	137	GLY	4.2
1	K	140	LYS	4.0
1	J	13	ASN	3.8
1	L	88	GLY	3.7
1	H	13	ASN	3.4

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Mol	Chain	Res	Type	RSRZ
1	B	12	THR	3.3
1	J	24	SER	3.3
1	E	12	THR	3.2
1	B	13	ASN	3.2
1	I	9	SER	3.1
1	A	13	ASN	3.0
1	I	10	LYS	2.9
1	I	140	LYS	2.9
1	L	13	ASN	2.8
1	L	105	LYS	2.8
1	L	15	LEU	2.7
1	L	24	SER	2.7
1	G	140	LYS	2.6
1	D	12	THR	2.6
1	J	14	LEU	2.5
1	G	13	ASN	2.5
1	G	27	LYS	2.4
1	L	110	ASP	2.4
1	K	137	GLY	2.3
1	H	100	SER	2.3
1	L	14	LEU	2.3
1	G	138	GLU	2.2
1	G	139	ALA	2.2
1	F	10	LYS	2.1
1	L	111	ILE	2.1
1	L	140	LYS	2.1
1	L	31	GLU	2.0
1	L	41	ILE	2.0
1	H	16	TYR	2.0
1	K	13	ASN	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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
2	TRS	G	607	8/8	0.85	0.23	5.77	49,52,53,56	0
2	TRS	F	606	8/8	0.92	0.19	3.04	45,50,52,52	0
2	TRS	C	603	8/8	0.91	0.18	2.37	32,35,39,41	0
2	TRS	J	612	8/8	0.90	0.19	1.77	64,66,68,68	0
2	TRS	K	611	8/8	0.93	0.26	1.40	58,63,64,64	0
2	TRS	F	605	8/8	0.85	0.26	1.01	65,66,66,67	0
2	TRS	E	601	8/8	0.97	0.11	-0.65	51,53,53,54	0
2	TRS	L	608	8/8	0.92	0.20	-1.62	81,82,82,84	0
2	TRS	A	602	8/8	0.70	0.33	-	83,83,84,85	0
2	TRS	K	610	8/8	0.60	0.33	-	86,89,91,92	0
2	TRS	I	609	8/8	0.59	0.35	-	124,125,126,126	0
2	TRS	C	604	8/8	0.67	0.28	-	77,79,79,80	0

6.5 Other polymers [i](#)

There are no such residues in this entry.