



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 9, 2016 – 06:21 PM EDT

PDB ID : 5IXM
Title : The LPS Transporter LptDE from Yersinia pestis, core complex
Authors : Botos, I.; Mayclin, S.J.; McCarthy, J.G.; Buchanan, S.K.
Deposited on : 2016-03-23
Resolution : 2.75 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20027674
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
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) : rb-20027674

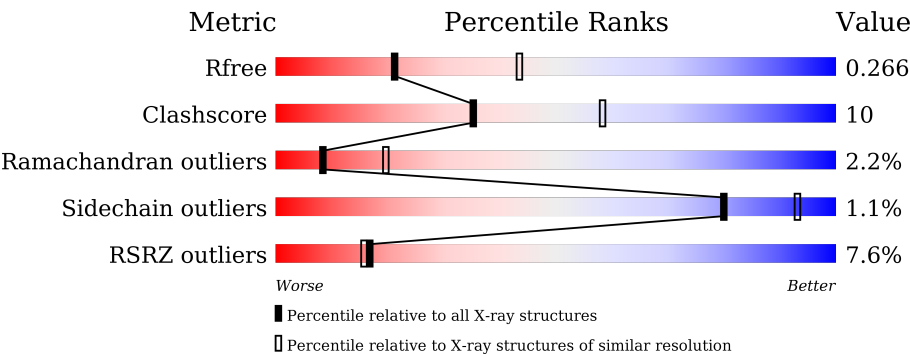
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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}	91344	3050 (2.78-2.70)
Clashscore	102246	3424 (2.78-2.70)
Ramachandran outliers	100387	3367 (2.78-2.70)
Sidechain outliers	100360	3368 (2.78-2.70)
RSRZ outliers	91569	3055 (2.78-2.70)

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	577	<div><div>4%</div><div><div></div><div>76%</div><div>17%</div><div>• 6%</div></div></div>
1	C	577	<div><div>5%</div><div><div></div><div>75%</div><div>18%</div><div>•• 6%</div></div></div>
1	E	577	<div><div>8%</div><div><div></div><div>72%</div><div>19%</div><div>• 7%</div></div></div>
1	G	577	<div><div>5%</div><div><div></div><div>74%</div><div>18%</div><div>• 7%</div></div></div>
2	B	198	<div><div>10%</div><div><div></div><div>54%</div><div>17%</div><div>• 29%</div></div></div>
2	D	198	<div><div>11%</div><div><div></div><div>58%</div><div>15%</div><div>• 26%</div></div></div>

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Mol	Chain	Length	Quality of chain
2	F	198	
2	H	198	

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
3	C8E	A	601	-	-	-	X
3	C8E	A	602	-	-	-	X
3	C8E	A	603	-	-	-	X
3	C8E	A	605	-	-	-	X
3	C8E	C	601	-	-	-	X
3	C8E	C	602	-	-	-	X
3	C8E	C	604	-	-	-	X
3	C8E	C	606	-	-	-	X
3	C8E	C	607	-	-	-	X
3	C8E	C	608	-	-	-	X
3	C8E	C	609	-	-	-	X
3	C8E	E	601	-	-	-	X
3	C8E	E	602	-	-	-	X
3	C8E	E	603	-	-	-	X
3	C8E	G	602	-	-	-	X
3	C8E	G	603	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 22831 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 LPS-assembly protein LptD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	543	Total	C	N	O	S	0	0	0
			4349	2765	742	829	13			
1	C	541	Total	C	N	O	S	0	0	0
			4346	2762	742	829	13			
1	E	538	Total	C	N	O	S	0	0	0
			4304	2737	736	818	13			
1	G	539	Total	C	N	O	S	0	0	0
			4329	2753	740	824	12			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	initiating methionine	UNP Q8ZIK3
A	-19	SER	-	expression tag	UNP Q8ZIK3
A	-18	ASN	-	expression tag	UNP Q8ZIK3
A	-17	HIS	-	expression tag	UNP Q8ZIK3
A	-16	HIS	-	expression tag	UNP Q8ZIK3
A	-15	HIS	-	expression tag	UNP Q8ZIK3
A	-14	HIS	-	expression tag	UNP Q8ZIK3
A	-13	HIS	-	expression tag	UNP Q8ZIK3
A	-12	HIS	-	expression tag	UNP Q8ZIK3
A	-11	HIS	-	expression tag	UNP Q8ZIK3
A	-10	HIS	-	expression tag	UNP Q8ZIK3
A	-9	HIS	-	expression tag	UNP Q8ZIK3
A	-8	HIS	-	expression tag	UNP Q8ZIK3
A	-7	GLU	-	expression tag	UNP Q8ZIK3
A	-6	ASN	-	expression tag	UNP Q8ZIK3
A	-5	LEU	-	expression tag	UNP Q8ZIK3
A	-4	TYR	-	expression tag	UNP Q8ZIK3
A	-3	PHE	-	expression tag	UNP Q8ZIK3
A	-2	GLN	-	expression tag	UNP Q8ZIK3
A	-1	SER	-	expression tag	UNP Q8ZIK3
A	0	MET	-	expression tag	UNP Q8ZIK3

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-20	MET	-	initiating methionine	UNP Q8ZIK3
C	-19	SER	-	expression tag	UNP Q8ZIK3
C	-18	ASN	-	expression tag	UNP Q8ZIK3
C	-17	HIS	-	expression tag	UNP Q8ZIK3
C	-16	HIS	-	expression tag	UNP Q8ZIK3
C	-15	HIS	-	expression tag	UNP Q8ZIK3
C	-14	HIS	-	expression tag	UNP Q8ZIK3
C	-13	HIS	-	expression tag	UNP Q8ZIK3
C	-12	HIS	-	expression tag	UNP Q8ZIK3
C	-11	HIS	-	expression tag	UNP Q8ZIK3
C	-10	HIS	-	expression tag	UNP Q8ZIK3
C	-9	HIS	-	expression tag	UNP Q8ZIK3
C	-8	HIS	-	expression tag	UNP Q8ZIK3
C	-7	GLU	-	expression tag	UNP Q8ZIK3
C	-6	ASN	-	expression tag	UNP Q8ZIK3
C	-5	LEU	-	expression tag	UNP Q8ZIK3
C	-4	TYR	-	expression tag	UNP Q8ZIK3
C	-3	PHE	-	expression tag	UNP Q8ZIK3
C	-2	GLN	-	expression tag	UNP Q8ZIK3
C	-1	SER	-	expression tag	UNP Q8ZIK3
C	0	MET	-	expression tag	UNP Q8ZIK3
E	-20	MET	-	initiating methionine	UNP Q8ZIK3
E	-19	SER	-	expression tag	UNP Q8ZIK3
E	-18	ASN	-	expression tag	UNP Q8ZIK3
E	-17	HIS	-	expression tag	UNP Q8ZIK3
E	-16	HIS	-	expression tag	UNP Q8ZIK3
E	-15	HIS	-	expression tag	UNP Q8ZIK3
E	-14	HIS	-	expression tag	UNP Q8ZIK3
E	-13	HIS	-	expression tag	UNP Q8ZIK3
E	-12	HIS	-	expression tag	UNP Q8ZIK3
E	-11	HIS	-	expression tag	UNP Q8ZIK3
E	-10	HIS	-	expression tag	UNP Q8ZIK3
E	-9	HIS	-	expression tag	UNP Q8ZIK3
E	-8	HIS	-	expression tag	UNP Q8ZIK3
E	-7	GLU	-	expression tag	UNP Q8ZIK3
E	-6	ASN	-	expression tag	UNP Q8ZIK3
E	-5	LEU	-	expression tag	UNP Q8ZIK3
E	-4	TYR	-	expression tag	UNP Q8ZIK3
E	-3	PHE	-	expression tag	UNP Q8ZIK3
E	-2	GLN	-	expression tag	UNP Q8ZIK3
E	-1	SER	-	expression tag	UNP Q8ZIK3
E	0	MET	-	expression tag	UNP Q8ZIK3

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-20	MET	-	initiating methionine	UNP Q8ZIK3
G	-19	SER	-	expression tag	UNP Q8ZIK3
G	-18	ASN	-	expression tag	UNP Q8ZIK3
G	-17	HIS	-	expression tag	UNP Q8ZIK3
G	-16	HIS	-	expression tag	UNP Q8ZIK3
G	-15	HIS	-	expression tag	UNP Q8ZIK3
G	-14	HIS	-	expression tag	UNP Q8ZIK3
G	-13	HIS	-	expression tag	UNP Q8ZIK3
G	-12	HIS	-	expression tag	UNP Q8ZIK3
G	-11	HIS	-	expression tag	UNP Q8ZIK3
G	-10	HIS	-	expression tag	UNP Q8ZIK3
G	-9	HIS	-	expression tag	UNP Q8ZIK3
G	-8	HIS	-	expression tag	UNP Q8ZIK3
G	-7	GLU	-	expression tag	UNP Q8ZIK3
G	-6	ASN	-	expression tag	UNP Q8ZIK3
G	-5	LEU	-	expression tag	UNP Q8ZIK3
G	-4	TYR	-	expression tag	UNP Q8ZIK3
G	-3	PHE	-	expression tag	UNP Q8ZIK3
G	-2	GLN	-	expression tag	UNP Q8ZIK3
G	-1	SER	-	expression tag	UNP Q8ZIK3
G	0	MET	-	expression tag	UNP Q8ZIK3

- Molecule 2 is a protein called LPS-assembly lipoprotein LptE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	141	Total	C	N	O	S	0	0	0
			1107	694	200	212	1			
2	D	146	Total	C	N	O	S	0	0	0
			1133	710	205	217	1			
2	F	145	Total	C	N	O	S	0	0	0
			1128	709	207	211	1			
2	H	147	Total	C	N	O	S	0	0	0
			1154	722	210	220	2			

There are 44 discrepancies between the modelled and reference sequences:

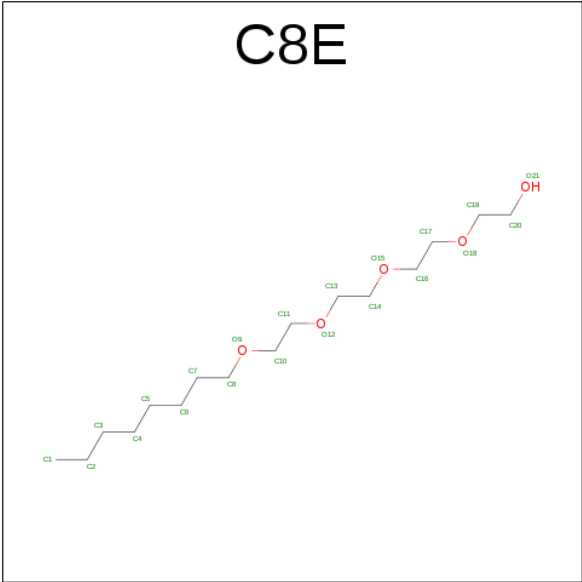
Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	ALA	-	expression tag	UNP Q7CJV2
B	-3	PRO	-	expression tag	UNP Q7CJV2
B	-2	ASN	-	expression tag	UNP Q7CJV2
B	-1	THR	-	expression tag	UNP Q7CJV2
B	0	SER	-	expression tag	UNP Q7CJV2

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Chain	Residue	Modelled	Actual	Comment	Reference
B	188	HIS	-	expression tag	UNP Q7CJV2
B	189	HIS	-	expression tag	UNP Q7CJV2
B	190	HIS	-	expression tag	UNP Q7CJV2
B	191	HIS	-	expression tag	UNP Q7CJV2
B	192	HIS	-	expression tag	UNP Q7CJV2
B	193	HIS	-	expression tag	UNP Q7CJV2
D	-4	ALA	-	expression tag	UNP Q7CJV2
D	-3	PRO	-	expression tag	UNP Q7CJV2
D	-2	ASN	-	expression tag	UNP Q7CJV2
D	-1	THR	-	expression tag	UNP Q7CJV2
D	0	SER	-	expression tag	UNP Q7CJV2
D	188	HIS	-	expression tag	UNP Q7CJV2
D	189	HIS	-	expression tag	UNP Q7CJV2
D	190	HIS	-	expression tag	UNP Q7CJV2
D	191	HIS	-	expression tag	UNP Q7CJV2
D	192	HIS	-	expression tag	UNP Q7CJV2
D	193	HIS	-	expression tag	UNP Q7CJV2
F	-4	ALA	-	expression tag	UNP Q7CJV2
F	-3	PRO	-	expression tag	UNP Q7CJV2
F	-2	ASN	-	expression tag	UNP Q7CJV2
F	-1	THR	-	expression tag	UNP Q7CJV2
F	0	SER	-	expression tag	UNP Q7CJV2
F	188	HIS	-	expression tag	UNP Q7CJV2
F	189	HIS	-	expression tag	UNP Q7CJV2
F	190	HIS	-	expression tag	UNP Q7CJV2
F	191	HIS	-	expression tag	UNP Q7CJV2
F	192	HIS	-	expression tag	UNP Q7CJV2
F	193	HIS	-	expression tag	UNP Q7CJV2
H	-4	ALA	-	expression tag	UNP Q7CJV2
H	-3	PRO	-	expression tag	UNP Q7CJV2
H	-2	ASN	-	expression tag	UNP Q7CJV2
H	-1	THR	-	expression tag	UNP Q7CJV2
H	0	SER	-	expression tag	UNP Q7CJV2
H	188	HIS	-	expression tag	UNP Q7CJV2
H	189	HIS	-	expression tag	UNP Q7CJV2
H	190	HIS	-	expression tag	UNP Q7CJV2
H	191	HIS	-	expression tag	UNP Q7CJV2
H	192	HIS	-	expression tag	UNP Q7CJV2
H	193	HIS	-	expression tag	UNP Q7CJV2

- Molecule 3 is (HYDROXYETHYLOXY)TRI(ETHYLOXY)OCTANE (three-letter code: C8E) (formula: C₁₆H₃₄O₅).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			21	16	5		
3	A	1	Total	C	O	0	0
			21	16	5		
3	A	1	Total	C	O	0	0
			21	16	5		
3	A	1	Total	C	O	0	0
			21	16	5		
3	A	1	Total	C	O	0	0
			21	16	5		
3	C	1	Total	C	O	0	0
			21	16	5		
3	C	1	Total	C	O	0	0
			21	16	5		
3	C	1	Total	C	O	0	0
			21	16	5		
3	C	1	Total	C	O	0	0
			21	16	5		
3	C	1	Total	C	O	0	0
			21	16	5		
3	C	1	Total	C	O	0	0
			21	16	5		
3	C	1	Total	C	O	0	0
			21	16	5		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	E	1	Total	C	O	0	0
			21	16	5		
3	E	1	Total	C	O	0	0
			21	16	5		
3	E	1	Total	C	O	0	0
			21	16	5		
3	E	1	Total	C	O	0	0
			21	16	5		
3	G	1	Total	C	O	0	0
			21	16	5		
3	G	1	Total	C	O	0	0
			21	16	5		
3	G	1	Total	C	O	0	0
			21	16	5		

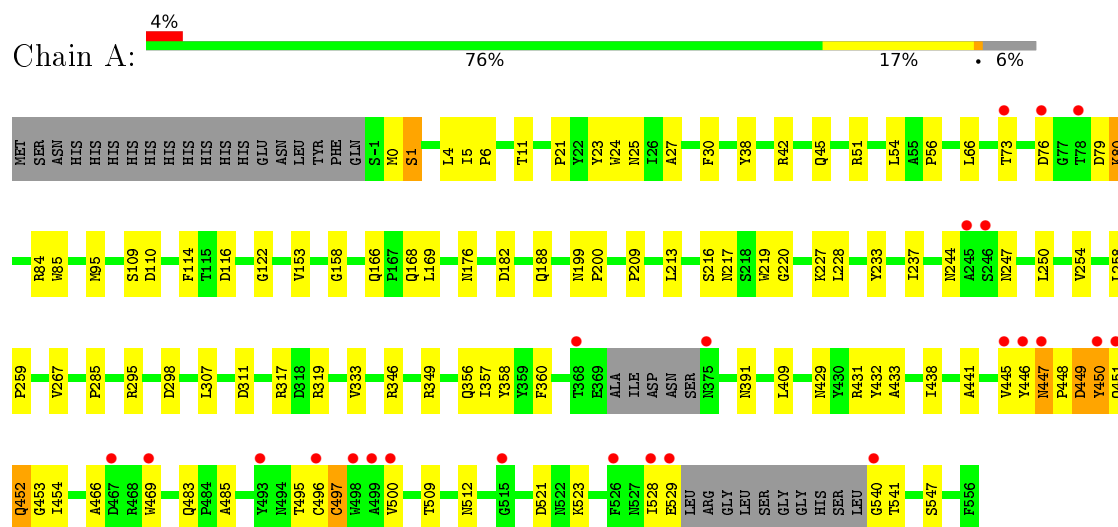
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	148	Total	O	0	0
			148	148		
4	B	38	Total	O	0	0
			38	38		
4	C	125	Total	O	0	0
			125	125		
4	D	27	Total	O	0	0
			27	27		
4	E	75	Total	O	0	0
			75	75		
4	F	16	Total	O	0	0
			16	16		
4	G	92	Total	O	0	0
			92	92		
4	H	19	Total	O	0	0
			19	19		

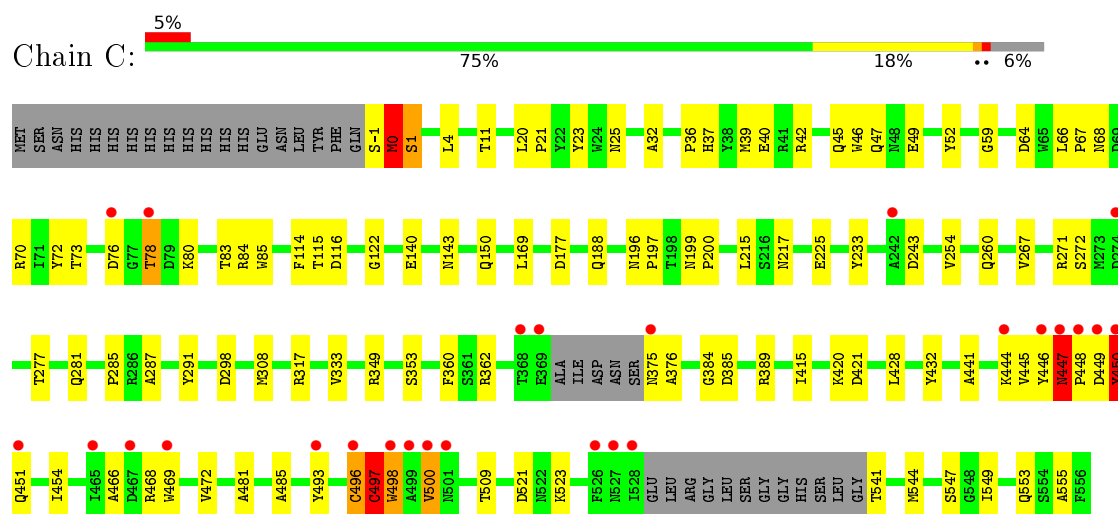
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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: LPS-assembly protein LptD

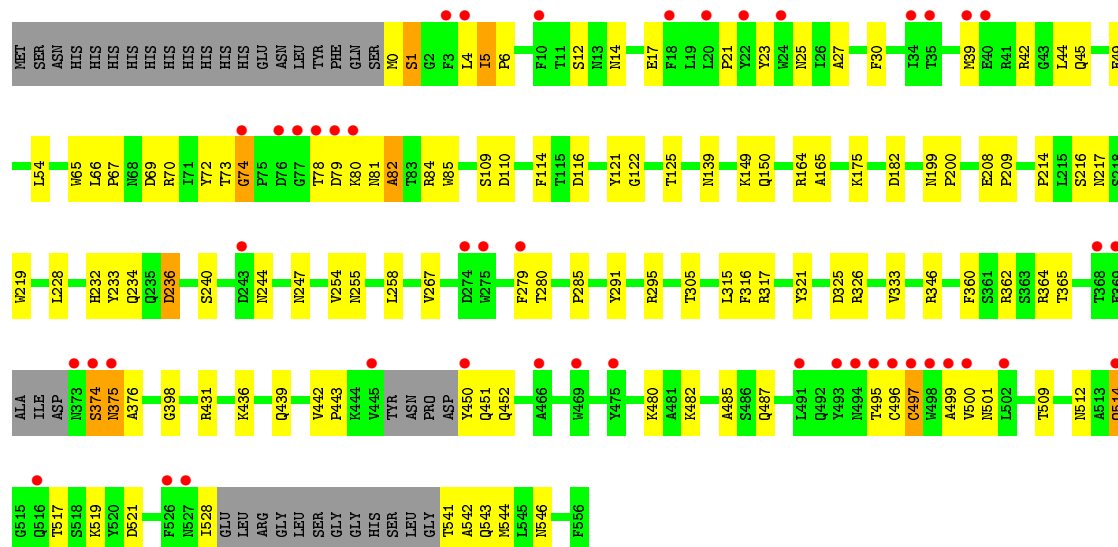


• Molecule 1: LPS-assembly protein LptD

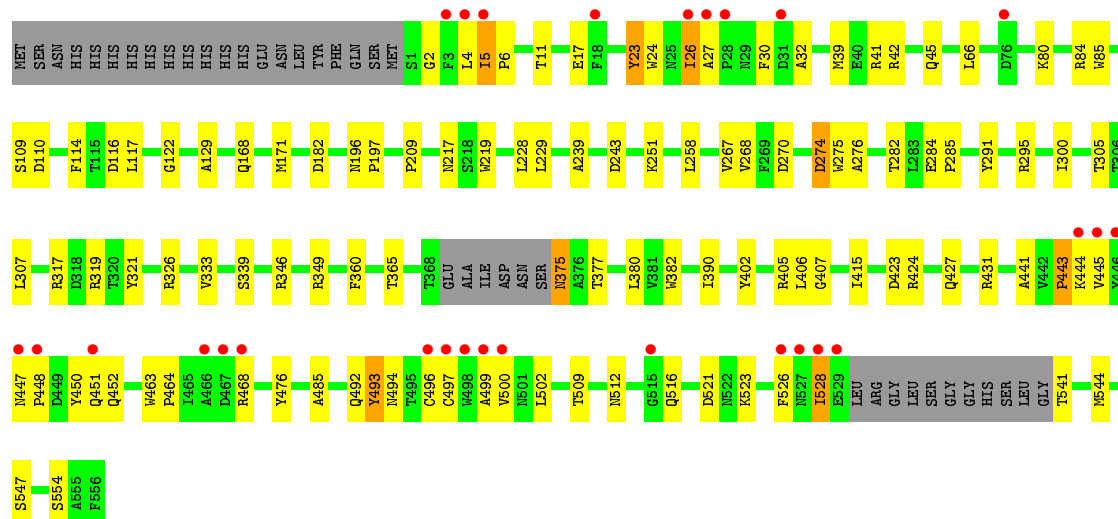
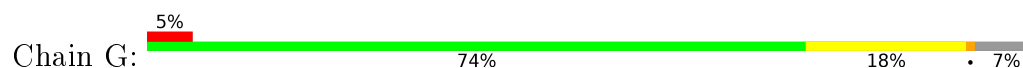


• Molecule 1: LPS-assembly protein LptD

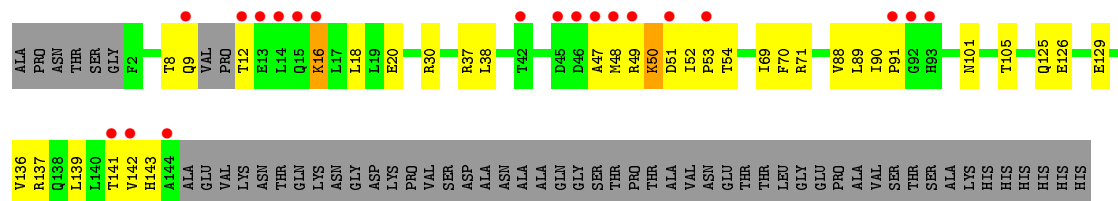




• Molecule 1: LPS-assembly protein LptD

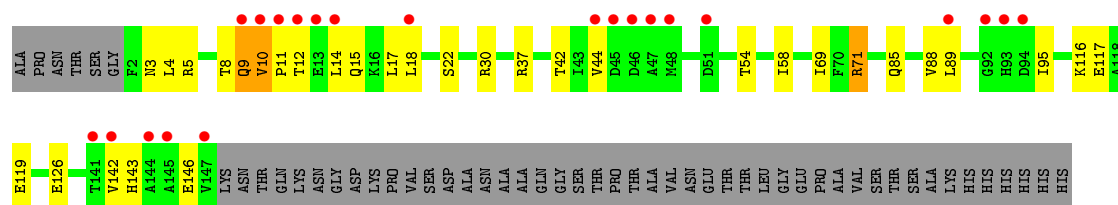


• Molecule 2: LPS-assembly lipoprotein LptE

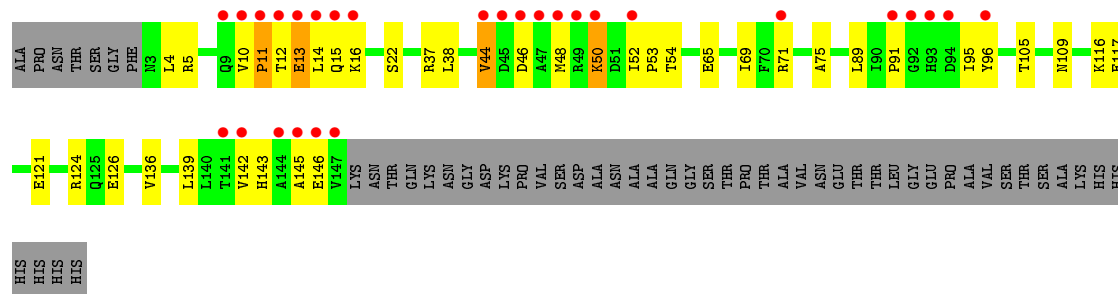


• Molecule 2: LPS-assembly lipoprotein LptE

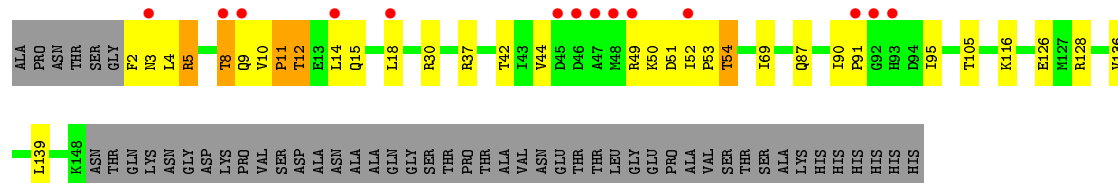




• Molecule 2: LPS-assembly lipoprotein LptE



• Molecule 2: LPS-assembly lipoprotein LptE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.89Å 176.35Å 143.85Å 90.00° 96.11° 90.00°	Depositor
Resolution (Å)	40.13 – 2.75 49.33 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.3 (40.13-2.75) 94.3 (49.33-2.75)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.88 (at 2.73Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.9_1692)	Depositor
R, R_{free}	0.212 , 0.261 0.215 , 0.266	Depositor DCC
R_{free} test set	5074 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	33.0	Xtriage
Anisotropy	0.450	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 63.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	22831	wwPDB-VP
Average B, all atoms (Å ²)	47.0	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 36.95 % 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 4.6183e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 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: C8E

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	0.26	0/4471	0.50	0/6091
1	C	0.27	0/4468	0.50	1/6087 (0.0%)
1	E	0.27	0/4423	0.52	0/6024
1	G	0.28	1/4451 (0.0%)	0.50	0/6065
2	B	0.28	0/1121	0.57	0/1523
2	D	0.29	0/1148	0.58	0/1564
2	F	0.29	0/1143	0.54	0/1555
2	H	0.26	0/1169	0.55	0/1589
All	All	0.27	1/22394 (0.0%)	0.52	1/30498 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	C	0	1
1	E	0	1
All	All	0	2

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	G	493	TYR	CD1-CE1	-5.17	1.31	1.39

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	447	ASN	N-CA-C	6.49	128.52	111.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	78	THR	Peptide
1	E	374	SER	Peptide

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	4349	0	4105	80	1
1	C	4346	0	4108	79	1
1	E	4304	0	4065	91	0
1	G	4329	0	4092	83	0
2	B	1107	0	1111	29	0
2	D	1133	0	1139	32	0
2	F	1128	0	1149	30	0
2	H	1154	0	1172	29	0
3	A	105	0	170	13	0
3	C	189	0	306	13	0
3	E	84	0	136	2	0
3	G	63	0	102	5	0
4	A	148	0	0	18	1
4	B	38	0	0	9	0
4	C	125	0	0	17	1
4	D	27	0	0	4	0
4	E	75	0	0	27	2
4	F	16	0	0	5	0
4	G	92	0	0	20	2
4	H	19	0	0	5	0
All	All	22831	0	21655	429	4

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 (429) 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:298:ASP:OD1	4:A:701:HOH:O	1.82	0.95
2:F:13:GLU:HG3	2:F:14:LEU:HG	1.51	0.93
2:B:12:THR:N	4:B:202:HOH:O	2.01	0.93
1:A:168:GLN:NE2	4:A:704:HOH:O	2.00	0.92
1:G:443:PRO:O	1:G:445:VAL:N	2.04	0.91
1:G:291:TYR:OH	4:G:701:HOH:O	1.90	0.88
1:E:232:HIS:ND1	4:E:707:HOH:O	2.09	0.86
1:E:482:LYS:HE3	2:H:8:THR:HG21	1.55	0.86
1:A:349:ARG:HH22	3:A:602:C8E:H42	1.41	0.85
1:A:429:ASN:OD1	4:A:702:HOH:O	1.94	0.85
1:G:168:GLN:NE2	4:G:706:HOH:O	2.09	0.85
1:A:54:LEU:O	4:A:703:HOH:O	1.94	0.84
1:G:502:LEU:HD13	1:G:526:PHE:HE1	1.42	0.84
1:E:39:MET:SD	4:E:768:HOH:O	2.33	0.84
1:G:447:ASN:HB3	1:G:450:TYR:HB3	1.61	0.83
1:G:349:ARG:NH1	4:G:708:HOH:O	2.12	0.82
1:C:497:CYS:SG	1:C:498:TRP:N	2.53	0.81
2:D:30:ARG:NH1	4:D:202:HOH:O	2.12	0.81
1:E:546:ASN:OD1	4:E:702:HOH:O	1.97	0.81
1:G:407:GLY:O	4:G:703:HOH:O	1.99	0.81
1:A:84:ARG:NH2	1:A:116:ASP:OD2	2.13	0.81
1:E:23:TYR:O	4:E:703:HOH:O	1.98	0.80
1:C:466:ALA:HB3	1:C:469:TRP:HB2	1.62	0.79
2:B:20:GLU:OE1	4:B:201:HOH:O	2.01	0.79
1:C:25:ASN:OD1	4:C:701:HOH:O	2.01	0.78
1:C:47:GLN:OE1	4:C:702:HOH:O	2.01	0.78
1:C:441:ALA:O	4:C:703:HOH:O	2.02	0.78
1:E:541:THR:O	1:E:543:GLN:N	2.17	0.78
1:A:176:ASN:OD1	4:A:705:HOH:O	2.01	0.77
1:E:72:TYR:O	1:E:74:GLY:N	2.18	0.77
1:G:4:LEU:HD11	1:G:23:TYR:HB3	1.66	0.77
1:E:79:ASP:OD2	1:E:82:ALA:HB2	1.84	0.76
1:E:139:ASN:O	4:E:704:HOH:O	2.03	0.76
2:D:22:SER:O	4:D:201:HOH:O	2.05	0.75
1:E:214:PRO:O	4:E:705:HOH:O	2.04	0.75
1:E:4:LEU:HD11	1:E:23:TYR:HB2	1.69	0.74
1:G:300:ILE:O	4:G:705:HOH:O	2.05	0.74
2:B:91:PRO:O	4:B:203:HOH:O	2.05	0.74
1:C:64:ASP:OD1	4:C:704:HOH:O	2.07	0.73
1:A:4:LEU:HB2	1:A:21:PRO:HG2	1.69	0.72
1:E:255:ASN:OD1	4:E:707:HOH:O	2.08	0.72
1:E:316:PHE:O	4:E:706:HOH:O	2.07	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:225:GLU:OE2	4:C:706:HOH:O	2.09	0.71
1:A:547:SER:O	4:A:706:HOH:O	2.09	0.71
1:G:24:TRP:HB3	1:G:32:ALA:HB3	1.72	0.71
1:G:502:LEU:HD13	1:G:526:PHE:CE1	2.25	0.71
1:E:315:LEU:O	4:E:709:HOH:O	2.09	0.70
1:G:468:ARG:HD2	1:G:493:TYR:HE1	1.57	0.70
2:F:14:LEU:HD12	2:F:53:PRO:HG2	1.73	0.70
2:B:16:LYS:NZ	2:B:52:ILE:HD12	2.07	0.70
2:F:116:LYS:NZ	4:F:203:HOH:O	2.25	0.69
1:A:521:ASP:OD2	4:A:707:HOH:O	2.09	0.69
1:E:279:PHE:O	4:E:708:HOH:O	2.09	0.69
1:A:158:GLY:O	4:A:708:HOH:O	2.11	0.69
1:E:69:ASP:N	1:E:81:ASN:O	2.24	0.69
1:E:70:ARG:N	4:E:701:HOH:O	1.86	0.69
1:E:501:ASN:HD22	1:E:528:ILE:HA	1.58	0.68
2:F:89:LEU:HG	2:F:95:ILE:HD12	1.74	0.68
1:G:415:ILE:HD11	1:G:427:GLN:HE21	1.57	0.68
2:H:87:GLN:HB3	2:H:95:ILE:HD11	1.75	0.67
1:G:84:ARG:NH2	1:G:116:ASP:OD2	2.26	0.67
1:C:84:ARG:NH2	1:C:116:ASP:OD2	2.24	0.67
2:B:125:GLN:OE1	4:B:204:HOH:O	2.12	0.66
1:A:483:GLN:OE1	4:A:710:HOH:O	2.14	0.66
2:B:16:LYS:HZ1	2:B:52:ILE:HD12	1.59	0.66
1:G:117:LEU:O	4:G:710:HOH:O	2.13	0.66
1:G:182:ASP:OD2	2:H:37:ARG:NH1	2.28	0.66
1:G:390:ILE:O	4:G:709:HOH:O	2.13	0.66
1:A:512:ASN:OD1	4:A:709:HOH:O	2.13	0.66
3:A:601:C8E:H141	1:C:384:GLY:HA3	1.76	0.66
1:E:480:LYS:O	4:E:711:HOH:O	2.13	0.65
1:E:234:GLN:OE1	4:E:710:HOH:O	2.13	0.65
1:C:11:THR:HG22	1:C:523:LYS:HG2	1.78	0.65
1:C:143:ASN:OD1	4:C:707:HOH:O	2.15	0.65
1:C:449:ASP:O	1:C:451:GLN:N	2.30	0.64
2:B:47:ALA:O	2:B:49:ARG:N	2.31	0.64
1:G:217:ASN:HB2	3:G:602:C8E:H171	1.80	0.64
1:E:495:THR:O	1:E:497:CYS:N	2.30	0.64
1:G:450:TYR:HD2	1:G:451:GLN:HG3	1.62	0.63
1:C:260:GLN:NE2	4:C:717:HOH:O	2.31	0.63
2:H:2:PHE:O	4:H:201:HOH:O	2.15	0.63
2:B:8:THR:HG22	2:B:9:GLN:H	1.65	0.62
2:D:126:GLU:OE1	4:D:203:HOH:O	2.16	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:26:ILE:HG22	4:G:712:HOH:O	1.98	0.62
2:B:88:VAL:HG11	2:B:142:VAL:HG11	1.82	0.62
2:B:30:ARG:NH1	4:B:207:HOH:O	2.22	0.61
1:A:259:PRO:HG3	3:C:602:C8E:H141	1.82	0.61
1:E:528:ILE:O	4:E:712:HOH:O	2.16	0.61
1:E:79:ASP:CG	1:E:82:ALA:HB2	2.20	0.61
1:C:298:ASP:O	4:C:708:HOH:O	2.16	0.61
1:G:448:PRO:O	1:G:452:GLN:HG3	2.00	0.61
2:D:10:VAL:O	2:D:12:THR:N	2.31	0.60
1:A:51:ARG:NH1	4:A:723:HOH:O	2.31	0.60
2:D:89:LEU:HG	2:D:95:ILE:HD12	1.81	0.60
1:G:468:ARG:HD2	1:G:493:TYR:CE1	2.34	0.60
1:E:244:ASN:HB3	1:E:247:ASN:HB2	1.82	0.60
1:A:1:SER:OG	1:A:24:TRP:O	2.11	0.60
2:D:9:GLN:O	2:D:10:VAL:HG23	2.02	0.60
1:C:272:SER:O	2:D:3:ASN:ND2	2.30	0.60
1:E:450:TYR:OH	2:F:71:ARG:NH1	2.34	0.60
3:A:603:C8E:H192	1:C:428:LEU:HD21	1.83	0.60
1:C:349:ARG:HH22	3:C:602:C8E:H191	1.67	0.59
1:C:496:CYS:O	1:C:497:CYS:HB3	2.01	0.59
1:G:4:LEU:HD11	1:G:23:TYR:CB	2.32	0.59
2:F:117:GLU:OE1	4:F:201:HOH:O	2.17	0.59
1:C:375:ASN:OD1	1:C:376:ALA:N	2.28	0.59
1:C:446:TYR:CD2	1:C:447:ASN:HB2	2.38	0.59
2:B:50:LYS:O	2:B:52:ILE:HG12	2.03	0.59
1:E:45:GLN:HG3	1:E:66:LEU:HB2	1.84	0.59
2:H:30:ARG:NH1	4:H:204:HOH:O	2.35	0.59
1:C:481:ALA:HB2	2:D:71:ARG:HH11	1.68	0.59
1:A:319:ARG:NH1	4:A:727:HOH:O	2.37	0.58
1:E:66:LEU:HD23	1:E:84:ARG:HG3	1.86	0.58
1:A:452:GLN:HG3	1:A:453:GLY:H	1.68	0.58
1:E:4:LEU:HB2	1:E:21:PRO:HG2	1.86	0.58
1:A:216:SER:OG	2:B:38:LEU:O	2.12	0.57
1:A:307:LEU:HD23	1:A:441:ALA:HA	1.84	0.57
1:A:495:THR:O	1:A:497:CYS:N	2.37	0.57
2:B:101:ASN:ND2	4:B:210:HOH:O	2.37	0.57
2:H:15:GLN:O	2:H:42:THR:N	2.36	0.57
1:A:466:ALA:HB3	1:A:469:TRP:HB2	1.87	0.56
1:C:285:PRO:HB2	3:C:601:C8E:H42	1.87	0.56
1:G:450:TYR:CD2	1:G:451:GLN:HG3	2.40	0.56
2:F:121:GLU:OE2	2:F:124:ARG:NH2	2.38	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:389:ARG:NH1	4:C:729:HOH:O	2.38	0.56
1:E:512:ASN:HB2	1:E:519:LYS:HE3	1.88	0.56
1:C:20:LEU:O	1:C:36:PRO:HD2	2.05	0.56
2:F:44:VAL:HG23	2:F:46:ASP:H	1.71	0.56
1:E:487:GLN:OE1	4:E:713:HOH:O	2.18	0.56
1:A:213:LEU:HD22	3:A:601:C8E:H61	1.87	0.55
1:C:73:THR:HB	1:C:78:THR:HB	1.89	0.55
2:B:129:GLU:OE2	4:B:205:HOH:O	2.18	0.55
1:G:485:ALA:HB1	2:H:69:ILE:HD11	1.89	0.55
1:G:492:GLN:NE2	1:G:494:ASN:HD21	2.04	0.55
1:E:326:ARG:HD2	1:E:365:THR:HB	1.89	0.55
1:A:485:ALA:HB1	2:B:69:ILE:HD11	1.88	0.55
2:D:10:VAL:C	2:D:12:THR:H	2.10	0.55
1:E:216:SER:OG	2:F:38:LEU:O	2.14	0.55
2:H:12:THR:HG23	2:H:14:LEU:H	1.72	0.55
2:H:52:ILE:O	4:H:203:HOH:O	2.17	0.55
1:A:244:ASN:HB3	1:A:247:ASN:HB2	1.90	0.54
1:E:499:ALA:O	4:E:714:HOH:O	2.18	0.54
1:A:540:GLY:N	4:A:731:HOH:O	2.41	0.54
1:A:311:ASP:OD2	4:A:711:HOH:O	2.17	0.54
1:A:446:TYR:OH	1:A:449:ASP:OD2	2.17	0.54
2:D:58:ILE:HD12	2:D:85:GLN:HG3	1.89	0.54
1:G:375:ASN:N	4:G:719:HOH:O	2.40	0.54
1:A:182:ASP:OD2	2:B:37:ARG:NH1	2.41	0.54
1:E:452:GLN:HE22	2:F:71:ARG:HG3	1.72	0.54
1:A:23:TYR:CE2	1:A:25:ASN:HB2	2.43	0.54
1:E:5:ILE:HD12	1:E:6:PRO:HD2	1.89	0.54
1:G:258:LEU:HD21	1:G:295:ARG:HB2	1.88	0.54
1:A:199:ASN:O	4:A:712:HOH:O	2.19	0.53
1:C:217:ASN:HB2	3:C:601:C8E:H161	1.90	0.53
2:H:4:LEU:HD23	2:H:5:ARG:H	1.73	0.53
1:A:391:ASN:HB2	3:A:602:C8E:H12	1.90	0.53
1:A:409:LEU:HD23	3:A:604:C8E:H192	1.91	0.53
1:C:4:LEU:HB2	1:C:21:PRO:HG2	1.89	0.53
1:C:233:TYR:HB2	1:C:254:VAL:HB	1.89	0.53
1:E:182:ASP:OD2	2:F:37:ARG:NH1	2.41	0.53
1:C:-1:SER:O	1:C:0:MET:HB2	2.10	0.53
1:E:175:LYS:NZ	4:E:734:HOH:O	2.42	0.53
1:E:69:ASP:HA	4:E:701:HOH:O	2.09	0.53
1:G:377:THR:HA	1:G:380:LEU:HD11	1.91	0.53
1:E:236:ASP:OD2	4:E:715:HOH:O	2.18	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:114:PHE:CE1	1:G:122:GLY:HA3	2.44	0.52
1:A:233:TYR:HB2	1:A:254:VAL:HB	1.90	0.52
1:A:349:ARG:NH2	3:A:602:C8E:H42	2.20	0.52
2:F:65:GLU:OE2	4:F:202:HOH:O	2.19	0.52
1:G:406:LEU:O	4:G:703:HOH:O	2.18	0.52
1:A:448:PRO:HA	1:A:451:GLN:O	2.09	0.52
1:E:66:LEU:HB3	1:E:84:ARG:HB2	1.91	0.51
1:A:333:VAL:HG23	1:A:360:PHE:HE2	1.75	0.51
1:A:346:ARG:O	2:B:141:THR:HG22	2.11	0.51
2:B:71:ARG:NE	4:B:206:HOH:O	2.22	0.51
1:C:150:GLN:NE2	4:C:728:HOH:O	2.38	0.51
1:E:431:ARG:NH2	2:F:105:THR:HG21	2.24	0.51
1:G:239:ALA:N	4:G:717:HOH:O	2.32	0.51
1:A:217:ASN:HB3	1:A:219:TRP:O	2.10	0.51
1:G:27:ALA:HB3	1:G:30:PHE:HB2	1.92	0.51
1:G:496:CYS:SG	1:G:497:CYS:N	2.81	0.51
1:C:493:TYR:HB3	1:C:500:VAL:HG22	1.92	0.51
1:G:415:ILE:HD11	1:G:427:GLN:NE2	2.25	0.51
1:E:0:MET:O	1:E:1:SER:HB2	2.11	0.51
1:A:432:TYR:HA	1:A:454:ILE:O	2.11	0.51
1:A:451:GLN:HE22	2:B:71:ARG:H	1.59	0.51
1:C:481:ALA:HB2	2:D:71:ARG:NH1	2.25	0.51
1:C:445:VAL:HG12	1:C:451:GLN:OE1	2.10	0.51
2:D:88:VAL:HG11	2:D:142:VAL:HG11	1.92	0.51
1:G:307:LEU:HD23	1:G:441:ALA:HA	1.92	0.51
1:A:11:THR:HG22	1:A:523:LYS:HG2	1.92	0.50
2:H:51:ASP:OD1	2:H:51:ASP:N	2.44	0.50
1:A:360:PHE:CE2	3:A:605:C8E:H171	2.46	0.50
1:A:448:PRO:O	1:A:450:TYR:N	2.45	0.50
1:C:333:VAL:HG23	1:C:360:PHE:HE2	1.77	0.50
1:A:509:THR:HG21	1:A:521:ASP:OD2	2.12	0.50
1:G:317:ARG:NH2	4:G:715:HOH:O	2.28	0.50
1:G:333:VAL:HG23	1:G:360:PHE:HE2	1.76	0.50
1:G:41:ARG:NH2	4:G:704:HOH:O	2.03	0.50
3:E:603:C8E:H21	3:G:602:C8E:H111	1.94	0.50
2:H:4:LEU:HD23	2:H:5:ARG:HG2	1.93	0.50
1:C:42:ARG:HG2	1:C:66:LEU:HD21	1.93	0.50
1:G:84:ARG:NH1	1:G:110:ASP:OD1	2.44	0.50
2:D:15:GLN:O	2:D:42:THR:N	2.41	0.50
2:H:44:VAL:HG21	2:H:52:ILE:HD11	1.94	0.50
2:B:137:ARG:NH2	4:B:209:HOH:O	2.32	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:8:THR:O	2:D:9:GLN:HB2	2.12	0.49
1:E:49:GLU:OE2	1:E:121:TYR:OH	2.31	0.49
2:D:14:LEU:HD11	2:D:17:LEU:HD23	1.94	0.49
1:E:398:GLY:HA3	3:E:603:C8E:H131	1.95	0.49
1:G:544:MET:O	1:G:547:SER:OG	2.31	0.49
1:G:11:THR:HG22	1:G:523:LYS:HG2	1.95	0.49
1:A:27:ALA:HB3	1:A:30:PHE:HD2	1.78	0.49
1:C:362:ARG:NH1	4:C:738:HOH:O	2.46	0.49
2:D:117:GLU:OE1	4:D:204:HOH:O	2.20	0.49
1:E:125:THR:HB	2:F:75:ALA:HB2	1.95	0.49
1:C:115:THR:O	4:C:711:HOH:O	2.20	0.49
1:E:17:GLU:HG2	1:E:39:MET:HG2	1.93	0.49
2:F:10:VAL:N	2:F:11:PRO:HD3	2.27	0.49
1:A:114:PHE:CE1	1:A:122:GLY:HA3	2.48	0.48
2:F:91:PRO:HB3	4:F:211:HOH:O	2.14	0.48
1:G:382:TRP:HB2	1:G:402:TYR:HB2	1.94	0.48
1:C:177:ASP:OD1	2:D:37:ARG:NH2	2.46	0.48
1:C:317:ARG:HD3	2:D:126:GLU:OE2	2.13	0.48
1:E:217:ASN:HB3	1:E:219:TRP:O	2.14	0.48
2:H:136:VAL:O	2:H:139:LEU:HB2	2.14	0.48
1:G:270:ASP:OD2	2:H:9:GLN:HB3	2.13	0.48
1:E:14:ASN:ND2	1:E:116:ASP:O	2.35	0.48
2:H:4:LEU:HD23	2:H:5:ARG:N	2.28	0.48
2:D:4:LEU:HD12	2:D:5:ARG:H	1.79	0.48
2:F:54:THR:OG1	2:F:89:LEU:HB2	2.13	0.48
2:H:3:ASN:O	4:H:201:HOH:O	2.20	0.48
3:A:603:C8E:H41	3:A:603:C8E:H13	1.73	0.47
1:G:554:SER:O	4:G:711:HOH:O	2.20	0.47
2:D:54:THR:OG1	2:D:89:LEU:HB2	2.14	0.47
1:E:436:LYS:NZ	4:E:738:HOH:O	2.47	0.47
1:A:356:GLN:HE22	1:A:358:TYR:HB2	1.78	0.47
1:C:509:THR:HG21	1:C:521:ASP:OD2	2.14	0.47
1:G:431:ARG:NH2	2:H:105:THR:HG21	2.30	0.47
2:H:10:VAL:N	2:H:11:PRO:HD3	2.29	0.47
1:G:45:GLN:HG3	1:G:66:LEU:HB2	1.96	0.47
3:C:603:C8E:H72	3:C:603:C8E:H41	1.73	0.47
1:E:442:VAL:HG12	2:F:109:ASN:HD22	1.79	0.47
1:A:227:LYS:HD3	1:A:319:ARG:HH22	1.79	0.47
1:C:23:TYR:CE2	1:C:25:ASN:HB2	2.49	0.47
1:A:267:VAL:O	1:A:285:PRO:HD2	2.15	0.47
1:A:45:GLN:HG3	1:A:66:LEU:HB2	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:56:PRO:HA	1:A:95:MET:HB2	1.97	0.47
1:C:544:MET:O	1:C:547:SER:OG	2.32	0.47
1:G:32:ALA:N	4:G:712:HOH:O	2.47	0.47
2:H:128:ARG:NH2	4:H:202:HOH:O	2.16	0.47
1:G:274:ASP:O	1:G:276:ALA:N	2.48	0.46
1:G:499:ALA:O	1:G:528:ILE:HG13	2.15	0.46
3:C:604:C8E:H52	3:C:604:C8E:H22	1.41	0.46
2:H:49:ARG:HH11	2:H:50:LYS:H	1.62	0.46
2:H:53:PRO:HB3	2:H:90:ILE:HG12	1.96	0.46
1:C:37:HIS:CD2	1:C:39:MET:HG3	2.51	0.46
3:C:602:C8E:H51	3:C:602:C8E:H22	1.41	0.46
2:H:116:LYS:HA	2:H:116:LYS:HD3	1.78	0.46
1:A:451:GLN:NE2	1:A:451:GLN:HA	2.30	0.46
2:H:3:ASN:OD1	2:H:4:LEU:N	2.49	0.46
1:A:438:ILE:HD12	1:A:453:GLY:HA2	1.98	0.46
1:E:199:ASN:O	4:E:717:HOH:O	2.21	0.46
1:G:423:ASP:OD1	1:G:424:ARG:HG3	2.15	0.46
1:A:166:GLN:HB3	4:A:751:HOH:O	2.16	0.46
2:D:8:THR:OG1	2:D:9:GLN:N	2.48	0.46
1:G:360:PHE:CG	3:G:603:C8E:H142	2.51	0.46
3:A:601:C8E:H72	3:A:601:C8E:H42	1.64	0.46
1:E:254:VAL:O	4:E:707:HOH:O	2.21	0.46
1:G:349:ARG:HH22	3:G:601:C8E:H62	1.81	0.46
1:G:42:ARG:HG2	1:G:66:LEU:HD21	1.96	0.46
1:C:485:ALA:HB1	2:D:69:ILE:HD11	1.98	0.45
1:E:236:ASP:OD1	1:E:236:ASP:N	2.44	0.45
1:A:317:ARG:HD3	2:B:126:GLU:OE2	2.16	0.45
1:E:150:GLN:HG3	1:E:164:ARG:NE	2.31	0.45
1:A:237:ILE:HD13	1:A:250:LEU:HB2	1.99	0.45
1:E:78:THR:HG22	1:E:79:ASP:N	2.31	0.45
1:A:209:PRO:HD2	1:A:228:LEU:O	2.17	0.45
1:C:353:SER:HB2	1:C:385:ASP:OD1	2.16	0.45
1:E:84:ARG:NH1	1:E:110:ASP:OD1	2.50	0.45
2:F:142:VAL:O	2:F:145:ALA:HB3	2.17	0.45
1:G:114:PHE:O	4:G:710:HOH:O	2.21	0.45
1:E:544:MET:N	4:E:730:HOH:O	2.48	0.45
1:C:150:GLN:HG2	4:C:728:HOH:O	2.16	0.45
1:C:0:MET:HB3	1:C:1:SER:H	1.35	0.45
1:E:233:TYR:HB2	1:E:254:VAL:HB	1.98	0.45
1:E:317:ARG:HD3	2:F:126:GLU:OE2	2.17	0.45
1:A:258:LEU:HD21	1:A:295:ARG:HB2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:602:C8E:H161	3:C:602:C8E:H191	1.69	0.45
1:C:36:PRO:HG3	1:C:46:TRP:CZ3	2.51	0.45
2:F:96:TYR:CD2	2:F:142:VAL:HG22	2.52	0.45
2:F:136:VAL:O	2:F:139:LEU:HB2	2.16	0.44
2:F:13:GLU:HG3	2:F:14:LEU:N	2.32	0.44
1:G:229:LEU:HD21	1:G:319:ARG:NH2	2.31	0.44
1:A:429:ASN:ND2	4:A:721:HOH:O	2.30	0.44
1:C:114:PHE:CE1	1:C:122:GLY:HA3	2.53	0.44
1:C:-1:SER:HB3	1:C:25:ASN:HD21	1.82	0.44
1:E:362:ARG:NH1	1:E:364:ARG:HD3	2.32	0.44
3:A:602:C8E:H191	3:A:602:C8E:H161	1.87	0.44
1:A:84:ARG:NH1	1:A:110:ASP:OD1	2.50	0.44
1:A:451:GLN:NE2	2:B:70:PHE:HB3	2.31	0.44
1:E:27:ALA:HB3	1:E:30:PHE:HB2	1.99	0.44
1:E:44:LEU:O	1:E:67:PRO:HD2	2.17	0.44
2:F:52:ILE:HA	2:F:53:PRO:HD3	1.87	0.44
1:G:317:ARG:NE	4:G:715:HOH:O	2.39	0.44
1:G:512:ASN:HB3	1:G:516:GLN:O	2.17	0.44
1:E:84:ARG:NH2	1:E:116:ASP:OD2	2.39	0.44
1:G:317:ARG:HD3	2:H:126:GLU:OE2	2.17	0.44
1:A:451:GLN:CD	2:B:70:PHE:HB3	2.38	0.44
1:E:78:THR:HG22	1:E:79:ASP:H	1.83	0.44
1:A:431:ARG:NH2	2:B:105:THR:HG21	2.33	0.44
1:A:449:ASP:C	1:A:451:GLN:H	2.21	0.44
1:A:73:THR:O	1:A:79:ASP:HB3	2.17	0.44
1:G:445:VAL:HG22	1:G:451:GLN:OE1	2.18	0.44
1:E:23:TYR:CE2	1:E:25:ASN:HB2	2.52	0.44
1:E:42:ARG:HD2	4:E:768:HOH:O	2.18	0.43
1:E:512:ASN:CG	1:E:514:GLN:HG3	2.37	0.43
1:G:380:LEU:HD13	1:G:405:ARG:HD3	2.00	0.43
1:C:72:TYR:O	1:C:73:THR:OG1	2.27	0.43
1:E:199:ASN:HA	1:E:200:PRO:HD3	1.86	0.43
1:A:333:VAL:HG23	1:A:360:PHE:CE2	2.53	0.43
1:C:169:LEU:O	1:C:188:GLN:HA	2.18	0.43
1:C:450:TYR:N	1:C:450:TYR:CD1	2.82	0.43
1:C:68:ASN:ND2	4:C:723:HOH:O	2.36	0.43
2:H:18:LEU:HD11	2:H:54:THR:HB	2.00	0.43
1:G:219:TRP:HB3	2:H:8:THR:HG22	2.00	0.43
1:C:466:ALA:CB	1:C:469:TRP:HB2	2.40	0.43
1:C:472:VAL:HG11	1:C:541:THR:HB	2.01	0.43
1:C:333:VAL:HG21	3:C:607:C8E:H132	2.01	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:500:VAL:HG13	4:A:722:HOH:O	2.19	0.43
3:C:605:C8E:H62	2:D:5:ARG:HG3	2.01	0.43
3:A:601:C8E:H41	3:A:601:C8E:H13	1.66	0.43
1:E:482:LYS:HD3	4:E:711:HOH:O	2.19	0.43
1:E:509:THR:HG21	1:E:521:ASP:OD2	2.18	0.43
1:G:509:THR:HG21	1:G:521:ASP:OD2	2.19	0.43
1:A:446:TYR:CD1	1:A:447:ASN:N	2.87	0.43
1:C:64:ASP:O	1:C:85:TRP:HA	2.19	0.43
2:F:15:GLN:N	4:F:210:HOH:O	2.51	0.43
1:A:356:GLN:HG2	1:A:357:ILE:N	2.34	0.43
1:A:85:TRP:CD1	1:A:109:SER:HB3	2.54	0.43
1:G:349:ARG:CZ	3:G:601:C8E:H102	2.49	0.43
2:D:143:HIS:O	2:D:146:GLU:HG2	2.19	0.42
1:C:277:THR:N	4:C:733:HOH:O	2.52	0.42
1:C:549:ILE:HD12	2:D:69:ILE:HD13	2.01	0.42
1:E:72:TYR:C	1:E:74:GLY:H	2.14	0.42
1:G:85:TRP:CD1	1:G:109:SER:HB3	2.54	0.42
1:E:209:PRO:HD2	1:E:228:LEU:O	2.19	0.42
2:H:90:ILE:HA	2:H:91:PRO:HD3	1.92	0.42
1:G:492:GLN:OE1	1:G:541:THR:N	2.53	0.42
1:C:215:LEU:HB3	3:C:601:C8E:H172	2.00	0.42
1:G:267:VAL:O	1:G:285:PRO:HD2	2.20	0.42
2:D:18:LEU:HD23	2:D:44:VAL:HB	2.02	0.42
1:G:282:THR:O	1:G:339:SER:HA	2.19	0.42
1:G:5:ILE:HG12	1:G:528:ILE:HG23	2.02	0.42
1:E:258:LEU:HD21	1:E:295:ARG:HB2	2.01	0.42
1:E:291:TYR:OH	4:E:716:HOH:O	2.20	0.42
1:E:30:PHE:CZ	1:E:54:LEU:HD21	2.55	0.42
1:E:72:TYR:HD2	1:E:79:ASP:OD2	2.02	0.42
1:G:463:TRP:HA	1:G:464:PRO:HD3	1.89	0.42
1:C:291:TYR:OH	4:C:712:HOH:O	2.21	0.42
1:C:70:ARG:NH1	4:C:736:HOH:O	2.45	0.42
2:F:143:HIS:O	2:F:146:GLU:HG2	2.19	0.42
1:G:209:PRO:HD2	1:G:228:LEU:O	2.20	0.42
1:A:169:LEU:O	1:A:188:GLN:HA	2.20	0.42
1:A:5:ILE:HA	1:A:6:PRO:HD3	1.84	0.42
2:B:136:VAL:O	2:B:139:LEU:HB2	2.20	0.42
1:C:32:ALA:HA	1:C:49:GLU:O	2.20	0.42
1:E:114:PHE:CE1	1:E:122:GLY:HA3	2.54	0.42
1:E:208:GLU:HA	1:E:228:LEU:O	2.19	0.42
1:E:485:ALA:HB1	2:F:69:ILE:HD11	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:129:ALA:HA	4:G:707:HOH:O	2.20	0.42
1:C:196:ASN:HA	1:C:197:PRO:HD3	1.87	0.41
1:C:287:ALA:HB2	3:C:601:C8E:H32	2.02	0.41
1:C:415:ILE:HD13	1:C:555:ALA:HB2	2.02	0.41
2:F:14:LEU:O	2:F:16:LYS:N	2.44	0.41
2:F:50:LYS:HB3	2:F:50:LYS:HE3	1.56	0.41
1:G:326:ARG:HD2	1:G:365:THR:HB	2.01	0.41
1:C:199:ASN:HA	1:C:200:PRO:HD3	1.87	0.41
1:C:52:TYR:CZ	1:C:59:GLY:HA3	2.55	0.41
1:G:251:LYS:HD2	1:G:251:LYS:HA	1.84	0.41
1:G:305:THR:HG23	1:G:321:TYR:HB3	2.03	0.41
2:H:8:THR:O	2:H:9:GLN:HB3	2.20	0.41
3:A:604:C8E:H13	3:A:604:C8E:H41	1.57	0.41
1:E:85:TRP:CD1	1:E:109:SER:HB3	2.56	0.41
1:G:346:ARG:HD3	1:G:346:ARG:HA	1.75	0.41
1:A:199:ASN:HA	1:A:200:PRO:HD3	1.87	0.41
1:C:45:GLN:HG3	1:C:66:LEU:HB2	2.03	0.41
1:C:67:PRO:HA	1:C:83:THR:HG23	2.03	0.41
1:E:149:LYS:HB2	1:E:165:ALA:HB3	2.02	0.41
1:E:280:THR:OG1	1:E:346:ARG:NH1	2.43	0.41
1:G:268:VAL:HG13	1:G:284:GLU:HG2	2.01	0.41
1:A:433:ALA:O	1:A:453:GLY:HA3	2.21	0.41
1:C:267:VAL:O	1:C:285:PRO:HD2	2.21	0.41
1:C:140:GLU:O	2:D:30:ARG:NH2	2.54	0.41
1:A:153:VAL:HG11	1:A:158:GLY:HA3	2.03	0.41
1:A:76:ASP:OD1	1:A:76:ASP:N	2.53	0.41
2:D:71:ARG:HH11	2:D:71:ARG:HD3	1.73	0.41
1:G:4:LEU:HD12	1:G:4:LEU:N	2.35	0.41
1:A:528:ILE:H	1:A:529:GLU:HG2	1.84	0.41
2:B:54:THR:OG1	2:B:89:LEU:HB2	2.21	0.41
2:B:90:ILE:HA	2:B:91:PRO:HD3	1.88	0.41
2:D:116:LYS:HD3	2:D:116:LYS:HA	1.78	0.41
1:E:333:VAL:HG23	1:E:360:PHE:HE2	1.86	0.41
1:E:375:ASN:CG	1:E:376:ALA:H	2.24	0.41
1:G:17:GLU:HG2	1:G:39:MET:HG2	2.02	0.41
2:B:53:PRO:HB3	2:B:90:ILE:HG12	2.02	0.41
1:E:267:VAL:O	1:E:285:PRO:HD2	2.21	0.41
1:E:436:LYS:HA	1:E:439:GLN:HG2	2.03	0.41
1:E:65:TRP:CZ3	1:E:67:PRO:HG3	2.56	0.41
2:F:11:PRO:HB2	2:F:12:THR:H	1.58	0.41
1:G:196:ASN:HA	1:G:197:PRO:HD3	1.94	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:271:ARG:NH2	1:C:281:GLN:OE1	2.54	0.40
1:C:420:LYS:HD3	1:C:421:ASP:OD1	2.19	0.40
1:E:495:THR:C	1:E:497:CYS:N	2.75	0.40
1:G:171:MET:SD	4:G:782:HOH:O	2.63	0.40
1:C:308:MET:SD	2:D:119:GLU:HG3	2.60	0.40
1:C:432:TYR:HA	1:C:454:ILE:O	2.21	0.40
1:G:476:TYR:OH	4:G:702:HOH:O	1.96	0.40
1:G:5:ILE:HA	1:G:6:PRO:HD3	1.92	0.40
1:C:45:GLN:OE1	1:C:47:GLN:NE2	2.40	0.40
2:D:14:LEU:CD1	2:D:17:LEU:HD23	2.52	0.40
2:D:9:GLN:O	2:D:10:VAL:CG2	2.68	0.40
1:E:233:TYR:OH	1:E:325:ASP:OD1	2.29	0.40
1:A:452:GLN:HG3	1:A:453:GLY:N	2.34	0.40
2:B:18:LEU:HB2	2:B:54:THR:HG22	2.04	0.40
1:A:42:ARG:HG2	1:A:66:LEU:HD21	2.02	0.40
1:C:349:ARG:NH2	3:C:602:C8E:H191	2.34	0.40
1:E:305:THR:HG23	1:E:321:TYR:HB3	2.04	0.40
1:E:70:ARG:NH2	4:E:718:HOH:O	2.21	0.40

All (4) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:A:846:HOH:O	4:C:823:HOH:O[2_746]	1.86	0.34
4:E:774:HOH:O	4:G:778:HOH:O[2_657]	2.12	0.08
4:E:774:HOH:O	4:G:780:HOH:O[2_657]	2.15	0.05
1:A:38:TYR:OH	1:C:40:GLU:OE2[2_846]	2.19	0.01

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	537/577 (93%)	490 (91%)	36 (7%)	11 (2%)	9	22
1	C	535/577 (93%)	491 (92%)	33 (6%)	11 (2%)	9	21
1	E	530/577 (92%)	483 (91%)	33 (6%)	14 (3%)	7	15
1	G	533/577 (92%)	488 (92%)	38 (7%)	7 (1%)	15	35
2	B	137/198 (69%)	123 (90%)	9 (7%)	5 (4%)	4	8
2	D	144/198 (73%)	134 (93%)	7 (5%)	3 (2%)	9	21
2	F	143/198 (72%)	131 (92%)	6 (4%)	6 (4%)	3	6
2	H	145/198 (73%)	135 (93%)	7 (5%)	3 (2%)	9	21
All	All	2704/3100 (87%)	2475 (92%)	169 (6%)	60 (2%)	8	20

All (60) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	1	SER
1	A	80	LYS
1	A	445	VAL
1	A	447	ASN
1	A	496	CYS
1	A	497	CYS
2	B	48	MET
2	B	51	ASP
1	C	1	SER
1	C	448	PRO
1	C	450	TYR
1	C	497	CYS
2	D	9	GLN
2	D	10	VAL
1	E	1	SER
1	E	73	THR
1	E	82	ALA
1	E	374	SER
1	E	496	CYS
1	E	497	CYS
1	E	500	VAL
1	E	514	GLN
1	E	542	ALA
2	F	5	ARG
2	F	11	PRO
2	F	13	GLU
2	F	44	VAL

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Mol	Chain	Res	Type
1	G	275	TRP
1	G	443	PRO
1	G	444	LYS
2	H	11	PRO
1	A	0	MET
1	C	0	MET
1	C	76	ASP
1	C	444	LYS
1	E	12	SER
2	F	4	LEU
2	F	48	MET
1	G	2	GLY
1	G	274	ASP
1	A	452	GLN
2	B	16	LYS
2	B	50	LYS
2	B	143	HIS
1	C	80	LYS
1	C	243	ASP
1	C	447	ASN
1	C	498	TRP
1	E	451	GLN
1	G	243	ASP
2	H	5	ARG
2	D	11	PRO
1	E	375	ASN
2	H	8	THR
1	A	449	ASP
1	A	541	THR
1	E	443	PRO
1	G	80	LYS
1	E	74	GLY
1	A	220	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	454/487 (93%)	452 (100%)	2 (0%)	93	98
1	C	456/487 (94%)	449 (98%)	7 (2%)	72	90
1	E	449/487 (92%)	444 (99%)	5 (1%)	80	93
1	G	453/487 (93%)	447 (99%)	6 (1%)	76	91
2	B	121/169 (72%)	121 (100%)	0	100	100
2	D	123/169 (73%)	122 (99%)	1 (1%)	86	95
2	F	122/169 (72%)	120 (98%)	2 (2%)	70	89
2	H	127/169 (75%)	125 (98%)	2 (2%)	70	89
All	All	2305/2624 (88%)	2280 (99%)	25 (1%)	80	93

All (25) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	80	LYS
1	A	450	TYR
1	C	0	MET
1	C	450	TYR
1	C	468	ARG
1	C	496	CYS
1	C	497	CYS
1	C	500	VAL
1	C	553	GLN
2	D	71	ARG
1	E	5	ILE
1	E	80	LYS
1	E	236	ASP
1	E	240	SER
1	E	517	THR
2	F	22	SER
2	F	50	LYS
1	G	5	ILE
1	G	23	TYR
1	G	26	ILE
1	G	375	ASN
1	G	500	VAL
1	G	528	ILE
2	H	12	THR
2	H	54	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (7) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	451	GLN
2	D	99	GLN
1	E	168	GLN
1	E	501	ASN
1	G	427	GLN
1	G	492	GLN
1	G	501	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 ⓘ

21 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	$\# Z > 2$
3	C8E	A	601	-	20,20,20	0.41	0	19,19,19	0.33	0
3	C8E	A	602	-	20,20,20	0.38	0	19,19,19	0.37	0
3	C8E	A	603	-	20,20,20	0.38	0	19,19,19	0.37	0
3	C8E	A	604	-	20,20,20	0.38	0	19,19,19	0.48	0
3	C8E	A	605	-	20,20,20	0.37	0	19,19,19	0.53	0
3	C8E	C	601	-	20,20,20	0.39	0	19,19,19	0.34	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	C8E	C	602	-	20,20,20	0.38	0	19,19,19	0.38	0
3	C8E	C	603	-	20,20,20	0.40	0	19,19,19	0.38	0
3	C8E	C	604	-	20,20,20	0.39	0	19,19,19	0.37	0
3	C8E	C	605	-	20,20,20	0.41	0	19,19,19	0.35	0
3	C8E	C	606	-	20,20,20	0.38	0	19,19,19	0.40	0
3	C8E	C	607	-	20,20,20	0.39	0	19,19,19	0.41	0
3	C8E	C	608	-	20,20,20	0.39	0	19,19,19	0.37	0
3	C8E	C	609	-	20,20,20	0.40	0	19,19,19	0.40	0
3	C8E	E	601	-	20,20,20	0.36	0	19,19,19	0.60	0
3	C8E	E	602	-	20,20,20	0.40	0	19,19,19	0.32	0
3	C8E	E	603	-	20,20,20	0.39	0	19,19,19	0.36	0
3	C8E	E	604	-	20,20,20	0.42	0	19,19,19	0.37	0
3	C8E	G	601	-	20,20,20	0.38	0	19,19,19	0.44	0
3	C8E	G	602	-	20,20,20	0.40	0	19,19,19	0.34	0
3	C8E	G	603	-	20,20,20	0.39	0	19,19,19	0.36	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	C8E	A	601	-	-	0/18/18/18	0/0/0/0
3	C8E	A	602	-	-	0/18/18/18	0/0/0/0
3	C8E	A	603	-	-	0/18/18/18	0/0/0/0
3	C8E	A	604	-	-	0/18/18/18	0/0/0/0
3	C8E	A	605	-	-	0/18/18/18	0/0/0/0
3	C8E	C	601	-	-	0/18/18/18	0/0/0/0
3	C8E	C	602	-	-	0/18/18/18	0/0/0/0
3	C8E	C	603	-	-	0/18/18/18	0/0/0/0
3	C8E	C	604	-	-	0/18/18/18	0/0/0/0
3	C8E	C	605	-	-	0/18/18/18	0/0/0/0
3	C8E	C	606	-	-	0/18/18/18	0/0/0/0
3	C8E	C	607	-	-	0/18/18/18	0/0/0/0
3	C8E	C	608	-	-	0/18/18/18	0/0/0/0
3	C8E	C	609	-	-	0/18/18/18	0/0/0/0
3	C8E	E	601	-	-	0/18/18/18	0/0/0/0
3	C8E	E	602	-	-	0/18/18/18	0/0/0/0
3	C8E	E	603	-	-	0/18/18/18	0/0/0/0
3	C8E	E	604	-	-	0/18/18/18	0/0/0/0
3	C8E	G	601	-	-	0/18/18/18	0/0/0/0
3	C8E	G	602	-	-	0/18/18/18	0/0/0/0
3	C8E	G	603	-	-	0/18/18/18	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

15 monomers are involved in 32 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	601	C8E	4	0
3	A	602	C8E	4	0
3	A	603	C8E	2	0
3	A	604	C8E	2	0
3	A	605	C8E	1	0
3	C	601	C8E	4	0
3	C	602	C8E	5	0
3	C	603	C8E	1	0
3	C	604	C8E	1	0
3	C	605	C8E	1	0
3	C	607	C8E	1	0
3	E	603	C8E	2	0
3	G	601	C8E	2	0
3	G	602	C8E	2	0
3	G	603	C8E	1	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	543/577 (94%)	0.06	24 (4%)	38 38	13, 32, 84, 131	0
1	C	541/577 (93%)	0.06	26 (4%)	34 34	11, 33, 89, 128	0
1	E	538/577 (93%)	0.50	45 (8%)	14 12	21, 53, 110, 166	0
1	G	539/577 (93%)	0.20	28 (5%)	31 31	18, 44, 97, 140	0
2	B	141/198 (71%)	0.55	20 (14%)	4 3	14, 36, 102, 127	0
2	D	146/198 (73%)	0.59	22 (15%)	3 2	13, 38, 106, 123	0
2	F	145/198 (73%)	0.84	28 (19%)	2 1	28, 54, 117, 145	0
2	H	147/198 (74%)	0.40	14 (9%)	10 9	19, 45, 101, 124	0
All	All	2740/3100 (88%)	0.29	207 (7%)	17 16	11, 41, 103, 166	0

All (207) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	500	VAL	9.1
2	F	11	PRO	8.6
1	A	446	TYR	7.9
2	B	142	VAL	6.5
2	F	14	LEU	6.4
1	A	245	ALA	6.3
2	B	48	MET	6.2
2	F	45	ASP	6.1
1	E	77	GLY	6.0
1	E	495	THR	6.0
2	F	48	MET	5.9
2	D	10	VAL	5.9
2	D	11	PRO	5.8
1	A	529	GLU	5.6
2	F	12	THR	5.5
2	D	48	MET	5.5

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Mol	Chain	Res	Type	RSRZ
2	B	141	THR	5.5
2	H	92	GLY	5.4
1	C	526	PHE	5.4
2	F	145	ALA	5.3
2	D	46	ASP	5.3
1	G	497	CYS	5.3
1	E	496	CYS	5.3
1	E	499	ALA	5.2
2	D	147	VAL	5.2
1	C	449	ASP	5.1
1	G	446	TYR	5.0
2	H	48	MET	5.0
2	D	12	THR	5.0
1	G	496	CYS	4.9
1	C	446	TYR	4.9
2	D	51	ASP	4.8
1	A	467	ASP	4.6
1	E	497	CYS	4.6
1	E	22	TYR	4.6
1	E	450	TYR	4.6
1	E	274	ASP	4.5
2	H	45	ASP	4.5
2	B	144	ALA	4.5
1	A	451	GLN	4.3
2	B	46	ASP	4.3
1	C	499	ALA	4.3
1	E	243	ASP	4.2
1	E	498	TRP	4.2
1	A	498	TRP	4.1
1	A	450	TYR	4.1
1	A	526	PHE	4.1
1	E	279	PHE	4.1
1	C	498	TRP	4.1
1	G	515	GLY	4.0
1	G	529	GLU	4.0
1	E	375	ASN	4.0
1	G	467	ASP	3.9
1	E	24	TRP	3.9
2	B	15	GLN	3.9
2	D	47	ALA	3.9
2	B	12	THR	3.9
2	H	49	ARG	3.9

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Mol	Chain	Res	Type	RSRZ
1	E	369	GLU	3.9
2	F	144	ALA	3.9
1	C	451	GLN	3.9
2	B	9	GLN	3.8
2	H	47	ALA	3.8
1	G	526	PHE	3.8
1	C	465	ILE	3.8
2	D	144	ALA	3.8
1	C	450	TYR	3.7
1	A	447	ASN	3.7
2	D	92	GLY	3.7
1	G	5	ILE	3.6
2	F	94	ASP	3.6
2	B	51	ASP	3.6
2	F	44	VAL	3.5
2	F	47	ALA	3.5
1	E	76	ASP	3.5
1	G	3	PHE	3.5
1	G	468	ARG	3.4
1	E	34	ILE	3.4
2	D	89	LEU	3.4
2	F	93	HIS	3.4
1	E	502	LEU	3.4
2	D	14	LEU	3.4
1	A	500	VAL	3.4
1	G	4	LEU	3.3
2	F	9	GLN	3.3
1	E	466	ALA	3.3
1	G	76	ASP	3.3
1	A	499	ALA	3.2
2	D	9	GLN	3.2
1	C	447	ASN	3.2
2	B	45	ASP	3.2
2	D	145	ALA	3.2
1	E	526	PHE	3.2
1	A	375	ASN	3.1
2	H	8	THR	3.1
2	F	10	VAL	3.1
1	A	73	THR	3.1
1	E	368	THR	3.1
2	F	91	PRO	3.1
2	B	49	ARG	3.1

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Mol	Chain	Res	Type	RSRZ
1	E	373	ASN	3.1
1	G	466	ALA	3.0
1	C	500	VAL	3.0
1	C	369	GLU	3.0
2	F	15	GLN	3.0
1	C	448	PRO	3.0
1	G	445	VAL	3.0
1	E	35	THR	3.0
1	C	496	CYS	3.0
1	G	498	TRP	3.0
1	E	74	GLY	3.0
1	G	451	GLN	2.9
1	G	528	ILE	2.9
1	A	528	ILE	2.9
1	E	79	ASP	2.9
2	D	142	VAL	2.9
2	F	49	ARG	2.9
1	E	78	THR	2.9
2	D	141	THR	2.9
2	B	91	PRO	2.9
2	D	13	GLU	2.9
2	B	47	ALA	2.9
2	H	52	ILE	2.9
1	E	469	TRP	2.8
1	E	80	LYS	2.8
1	G	26	ILE	2.8
1	E	494	ASN	2.8
2	F	52	ILE	2.8
1	G	444	LYS	2.8
1	E	275	TRP	2.8
2	D	45	ASP	2.8
2	D	94	ASP	2.7
1	E	493	TYR	2.7
2	F	96	TYR	2.7
2	F	46	ASP	2.7
1	G	28	PRO	2.7
1	A	496	CYS	2.7
2	F	146	GLU	2.7
1	G	448	PRO	2.7
2	F	50	LYS	2.6
1	A	493	TYR	2.6
2	B	92	GLY	2.6

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Mol	Chain	Res	Type	RSRZ
2	H	91	PRO	2.6
1	G	500	VAL	2.6
1	C	527	ASN	2.6
1	A	540	GLY	2.6
1	C	528	ILE	2.6
2	F	142	VAL	2.6
1	C	274	ASP	2.6
2	B	13	GLU	2.6
2	H	14	LEU	2.6
2	F	147	VAL	2.5
1	C	444	LYS	2.5
1	E	4	LEU	2.5
1	E	514	GLN	2.5
1	A	368	THR	2.5
2	H	46	ASP	2.5
1	A	445	VAL	2.5
1	E	445	VAL	2.5
2	B	42	THR	2.4
2	F	141	THR	2.4
1	C	375	ASN	2.4
1	C	242	ALA	2.4
2	F	16	LYS	2.4
1	E	491	LEU	2.4
1	G	27	ALA	2.4
1	E	374	SER	2.4
2	D	18	LEU	2.4
1	C	493	TYR	2.3
2	H	18	LEU	2.3
1	A	76	ASP	2.3
2	B	93	HIS	2.3
2	H	3	ASN	2.3
2	H	9	GLN	2.3
1	G	31	ASP	2.3
1	G	499	ALA	2.2
1	A	515	GLY	2.2
1	G	18	PHE	2.2
1	E	40	GLU	2.2
2	F	71	ARG	2.2
1	E	475	TYR	2.2
1	E	527	ASN	2.2
1	G	447	ASN	2.2
1	E	3	PHE	2.2

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Mol	Chain	Res	Type	RSRZ
1	E	39	MET	2.2
1	C	78	THR	2.2
1	C	469	TRP	2.2
1	E	10	PHE	2.2
2	F	92	GLY	2.1
1	E	18	PHE	2.1
1	E	516	GLN	2.1
2	D	44	VAL	2.1
2	B	14	LEU	2.1
2	H	93	HIS	2.1
2	D	93	HIS	2.1
1	G	527	ASN	2.1
2	B	16	LYS	2.1
2	B	53	PRO	2.1
2	F	13	GLU	2.1
1	C	501	ASN	2.1
1	A	78	THR	2.1
1	C	467	ASP	2.1
1	A	246	SER	2.1
1	C	368	THR	2.0
1	C	76	ASP	2.0
1	A	469	TRP	2.0
1	E	20	LEU	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [i](#)

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

6.3 Carbohydrates [i](#)

There are no 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(\AA^2)	Q<0.9
3	C8E	G	603	21/21	0.87	0.43	7.17	33,76,105,109	0
3	C8E	C	608	21/21	0.82	0.29	6.18	31,56,78,86	0
3	C8E	C	604	21/21	0.83	0.41	5.94	46,65,83,90	0
3	C8E	E	603	21/21	0.85	0.30	5.35	20,47,69,75	0
3	C8E	E	602	21/21	0.90	0.34	5.10	24,42,69,82	0
3	C8E	C	607	21/21	0.88	0.29	4.99	26,54,74,94	0
3	C8E	A	601	21/21	0.85	0.28	4.81	24,38,60,74	0
3	C8E	C	602	21/21	0.88	0.25	4.81	30,47,60,73	0
3	C8E	A	603	21/21	0.86	0.36	4.14	31,65,88,99	0
3	C8E	E	601	21/21	0.92	0.24	4.06	33,48,86,89	0
3	C8E	A	605	21/21	0.90	0.31	3.61	31,57,109,114	0
3	C8E	C	601	21/21	0.92	0.27	3.60	29,44,80,96	0
3	C8E	G	602	21/21	0.89	0.31	3.46	35,55,115,123	0
3	C8E	C	606	21/21	0.90	0.26	3.30	22,48,69,71	0
3	C8E	C	609	21/21	0.89	0.24	2.54	21,60,103,110	0
3	C8E	A	602	21/21	0.89	0.23	2.53	28,44,65,72	0
3	C8E	E	604	21/21	0.84	0.32	1.77	38,63,85,93	0
3	C8E	C	603	21/21	0.94	0.23	1.59	25,47,71,74	0
3	C8E	G	601	21/21	0.93	0.22	1.23	21,48,62,73	0
3	C8E	A	604	21/21	0.94	0.20	1.16	22,34,46,49	0
3	C8E	C	605	21/21	0.88	0.23	1.02	38,52,71,88	0

6.5 Other polymers ⓘ

There are no such residues in this entry.