



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 9, 2016 – 03:07 PM EDT

PDB ID : 5IV8
Title : The LPS Transporter LptDE from *Klebsiella pneumoniae*, core complex
Authors : Botos, I.; McCarthy, J.G.; Buchanan, S.K.
Deposited on : 2016-03-20
Resolution : 2.94 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

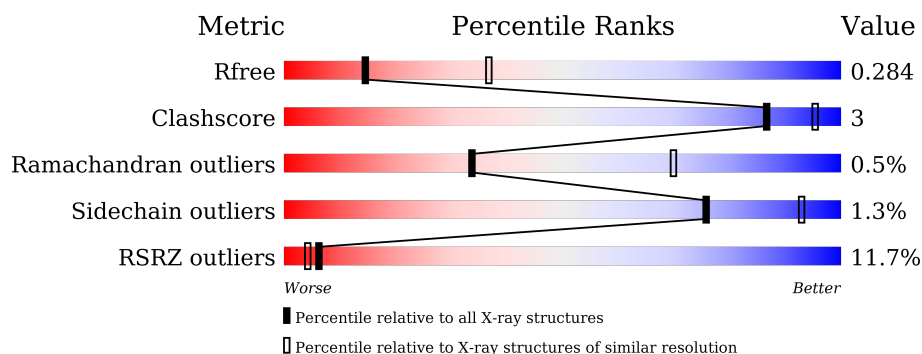
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.94 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	2057 (2.98-2.90)
Clashscore	102246	2308 (2.98-2.90)
Ramachandran outliers	100387	2245 (2.98-2.90)
Sidechain outliers	100360	2247 (2.98-2.90)
RSRZ outliers	91569	2065 (2.98-2.90)

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	601	<div> <div>7%</div> <div>84%</div> <div>8%</div> <div>8%</div> </div>
1	C	601	<div> <div>17%</div> <div>80%</div> <div>10%</div> <div>10%</div> </div>
2	B	182	<div> <div>7%</div> <div>75%</div> <div>5%</div> <div>20%</div> </div>
2	D	182	<div> <div>5%</div> <div>73%</div> <div>9%</div> <div>18%</div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	C8E	A	801	-	-	-	X
3	C8E	A	802	-	-	-	X
3	C8E	C	801	-	-	-	X
3	C8E	C	802	-	-	-	X
3	C8E	C	803	-	-	-	X
3	C8E	C	804	-	-	-	X
3	C8E	D	201	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 11347 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 biosynthesis protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	552	Total	C	N	O	S	0	0	0
			4472	2808	766	888	10			
1	C	543	Total	C	N	O	S	0	0	0
			4401	2767	749	875	10			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	181	MET	-	initiating methionine	UNP A0A0U3IWD2
A	182	SER	-	expression tag	UNP A0A0U3IWD2
A	183	ASN	-	expression tag	UNP A0A0U3IWD2
A	184	HIS	-	expression tag	UNP A0A0U3IWD2
A	185	HIS	-	expression tag	UNP A0A0U3IWD2
A	186	HIS	-	expression tag	UNP A0A0U3IWD2
A	187	HIS	-	expression tag	UNP A0A0U3IWD2
A	188	HIS	-	expression tag	UNP A0A0U3IWD2
A	189	HIS	-	expression tag	UNP A0A0U3IWD2
A	190	HIS	-	expression tag	UNP A0A0U3IWD2
A	191	HIS	-	expression tag	UNP A0A0U3IWD2
A	192	HIS	-	expression tag	UNP A0A0U3IWD2
A	193	HIS	-	expression tag	UNP A0A0U3IWD2
A	194	GLU	-	expression tag	UNP A0A0U3IWD2
A	195	ASN	-	expression tag	UNP A0A0U3IWD2
A	196	LEU	-	expression tag	UNP A0A0U3IWD2
A	197	TYR	-	expression tag	UNP A0A0U3IWD2
A	198	PHE	-	expression tag	UNP A0A0U3IWD2
A	199	GLN	-	expression tag	UNP A0A0U3IWD2
A	200	SER	-	expression tag	UNP A0A0U3IWD2
A	201	MET	-	expression tag	UNP A0A0U3IWD2
C	181	MET	-	initiating methionine	UNP A0A0U3IWD2
C	182	SER	-	expression tag	UNP A0A0U3IWD2
C	183	ASN	-	expression tag	UNP A0A0U3IWD2
C	184	HIS	-	expression tag	UNP A0A0U3IWD2

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Chain	Residue	Modelled	Actual	Comment	Reference
C	185	HIS	-	expression tag	UNP A0A0U3IWD2
C	186	HIS	-	expression tag	UNP A0A0U3IWD2
C	187	HIS	-	expression tag	UNP A0A0U3IWD2
C	188	HIS	-	expression tag	UNP A0A0U3IWD2
C	189	HIS	-	expression tag	UNP A0A0U3IWD2
C	190	HIS	-	expression tag	UNP A0A0U3IWD2
C	191	HIS	-	expression tag	UNP A0A0U3IWD2
C	192	HIS	-	expression tag	UNP A0A0U3IWD2
C	193	HIS	-	expression tag	UNP A0A0U3IWD2
C	194	GLU	-	expression tag	UNP A0A0U3IWD2
C	195	ASN	-	expression tag	UNP A0A0U3IWD2
C	196	LEU	-	expression tag	UNP A0A0U3IWD2
C	197	TYR	-	expression tag	UNP A0A0U3IWD2
C	198	PHE	-	expression tag	UNP A0A0U3IWD2
C	199	GLN	-	expression tag	UNP A0A0U3IWD2
C	200	SER	-	expression tag	UNP A0A0U3IWD2
C	201	MET	-	expression tag	UNP A0A0U3IWD2

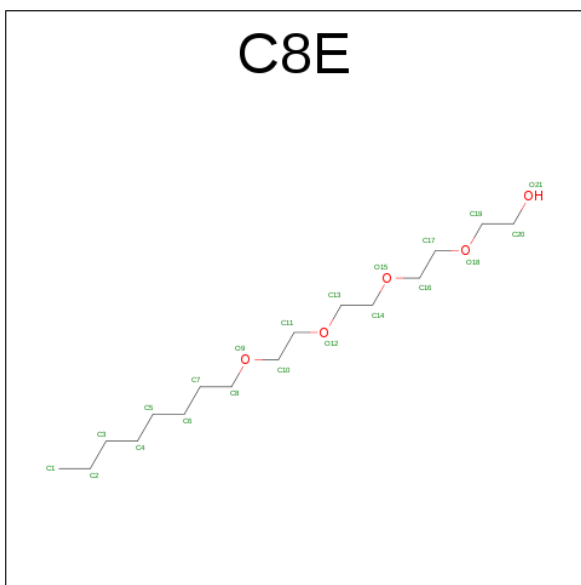
- Molecule 2 is a protein called LPS-assembly lipoprotein LptE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	146	Total	C	N	O	S	0	0	0
			1140	713	201	220	6			
2	D	149	Total	C	N	O	S	0	0	0
			1164	727	205	226	6			

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	15	ALA	-	expression tag	UNP A0A0J4W1Y0
B	16	PRO	-	expression tag	UNP A0A0J4W1Y0
B	17	ASN	-	expression tag	UNP A0A0J4W1Y0
B	18	THR	-	expression tag	UNP A0A0J4W1Y0
B	19	SER	-	expression tag	UNP A0A0J4W1Y0
D	15	ALA	-	expression tag	UNP A0A0J4W1Y0
D	16	PRO	-	expression tag	UNP A0A0J4W1Y0
D	17	ASN	-	expression tag	UNP A0A0J4W1Y0
D	18	THR	-	expression tag	UNP A0A0J4W1Y0
D	19	SER	-	expression tag	UNP A0A0J4W1Y0

- 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	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	D	1	Total	C	O	0	0
			21	16	5		

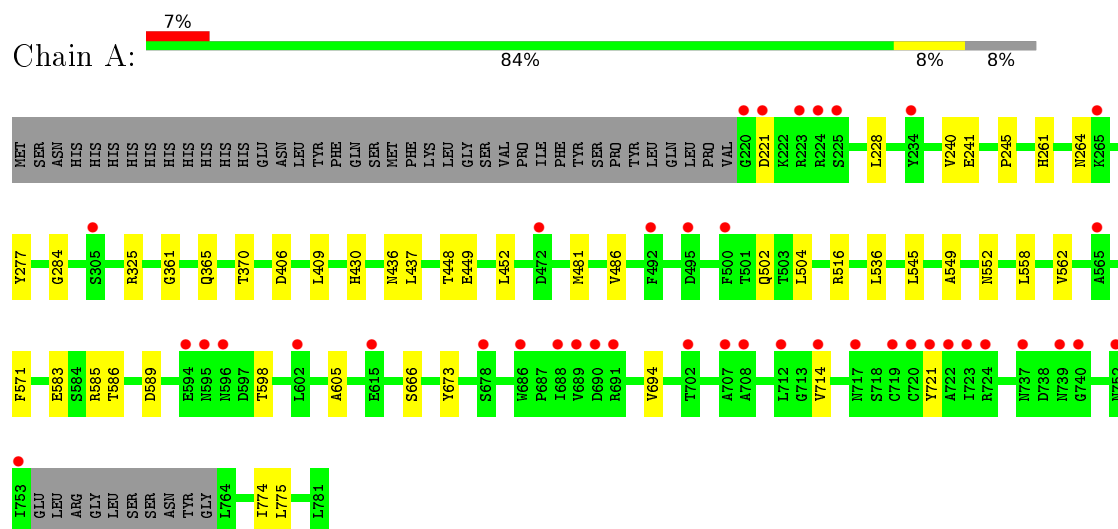
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	2	Total	O	0	0
			2	2		

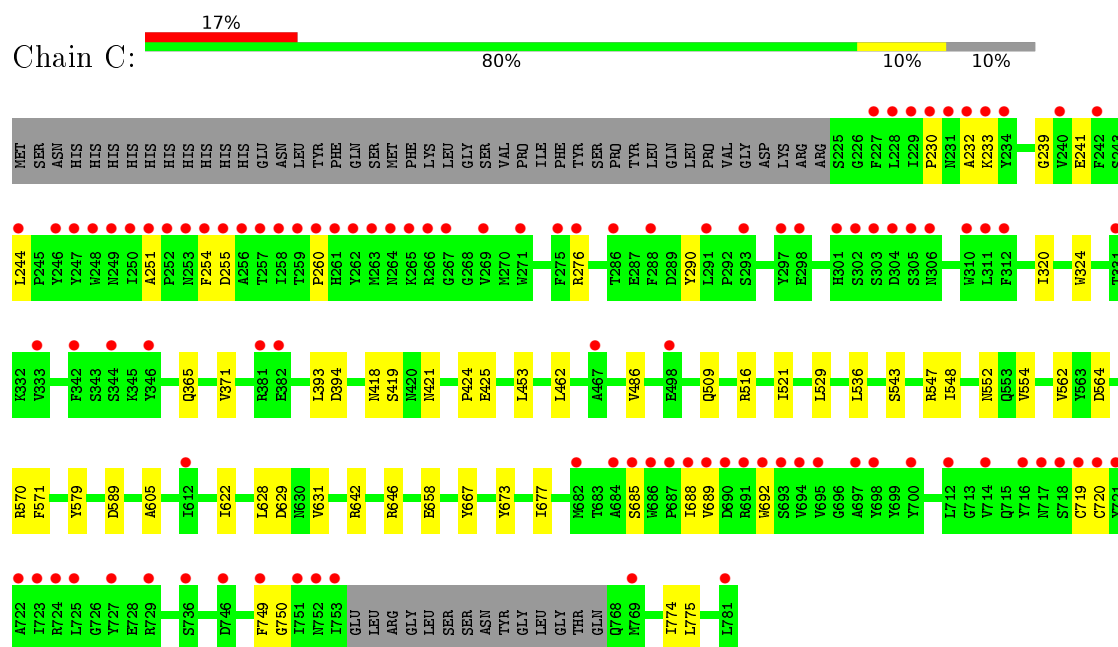
3 Residue-property plots

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.

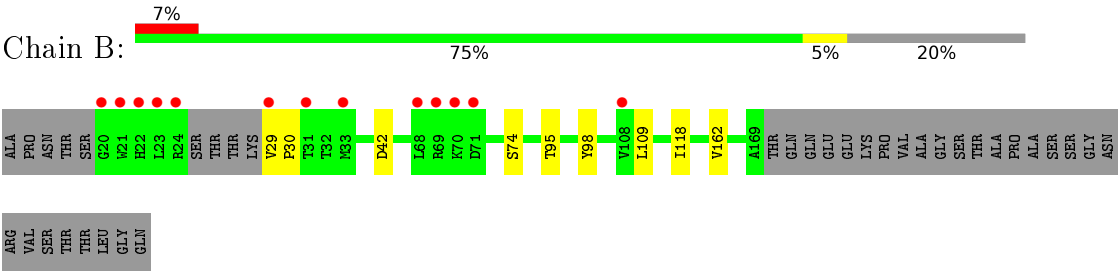
- Molecule 1: LPS biosynthesis protein



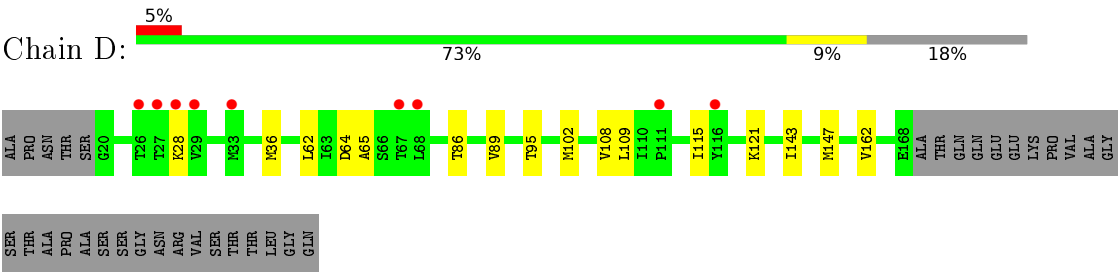
- Molecule 1: LPS biosynthesis protein



- 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	75.02Å 173.05Å 84.76Å 90.00° 111.26° 90.00°	Depositor
Resolution (Å)	41.41 – 2.94 46.58 – 2.94	Depositor EDS
% Data completeness (in resolution range)	83.7 (41.41-2.94) 83.7 (46.58-2.94)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.03 (at 2.96Å)	Xtriage
Refinement program	PHENIX (1.10 _2155: ???)	Depositor
R, R_{free}	0.235 , 0.284 0.230 , 0.284	Depositor DCC
R_{free} test set	1987 reflections (5.54%)	DCC
Wilson B-factor (Å ²)	49.0	Xtriage
Anisotropy	0.284	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 49.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.88	EDS
Total number of atoms	11347	wwPDB-VP
Average B, all atoms (Å ²)	72.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.69% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: 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	0.24	0/4588	0.43	0/6240
1	C	0.24	0/4517	0.45	0/6147
2	B	0.24	0/1158	0.41	0/1569
2	D	0.24	0/1183	0.41	0/1604
All	All	0.24	0/11446	0.43	0/15560

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4472	0	4149	21	0
1	C	4401	0	4074	27	0
2	B	1140	0	1147	5	0
2	D	1164	0	1175	8	0
3	A	63	0	102	2	0
3	C	84	0	136	1	0
3	D	21	0	34	0	0
4	A	2	0	0	0	0
All	All	11347	0	10817	60	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 3.

All (60) 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:C:642:ARG:HE	1:C:646:ARG:HD3	1.52	0.74
1:C:646:ARG:HA	1:C:685:SER:O	1.91	0.70
1:C:251:ALA:HB3	1:C:254:PHE:HB2	1.78	0.64
1:A:774:ILE:HG13	1:A:775:LEU:HD12	1.82	0.62
2:D:109:LEU:HG	2:D:115:ILE:HG22	1.83	0.60
1:C:453:LEU:HD13	1:C:548:ILE:HD13	1.85	0.58
1:C:320:ILE:HG13	1:C:324:TRP:HB2	1.87	0.56
1:C:233:LYS:HD2	1:C:750:GLY:HA3	1.88	0.56
1:A:452:LEU:HB2	3:A:801:C8E:H62	1.89	0.54
1:A:409:LEU:HB2	3:A:801:C8E:H171	1.89	0.54
1:C:775:LEU:HD21	2:D:89:VAL:HG13	1.89	0.53
2:D:108:VAL:HG11	2:D:162:VAL:HG21	1.90	0.53
1:C:562:VAL:HB	1:C:571:PHE:HB3	1.91	0.53
2:B:118:ILE:HD11	2:B:162:VAL:HG21	1.92	0.52
1:C:529:LEU:HD21	1:C:543:SER:HB3	1.92	0.51
1:A:361:GLY:HA3	1:A:370:THR:HG22	1.92	0.51
1:A:536:LEU:HD11	1:A:605:ALA:HB3	1.93	0.51
1:C:230:PRO:HB3	1:C:244:LEU:HD23	1.91	0.51
2:D:102:MET:O	2:D:121:LYS:HA	2.11	0.50
1:C:255:ASP:HB3	1:C:276:ARG:HB2	1.93	0.49
1:A:562:VAL:HG22	1:A:571:PHE:HB3	1.95	0.49
2:B:74:SER:HB3	2:B:109:LEU:HB2	1.94	0.48
1:A:277:TYR:CZ	1:A:284:GLY:HA3	2.49	0.47
1:A:448:THR:HG22	1:A:486:VAL:HG12	1.96	0.47
1:A:325:ARG:NH2	2:B:42:ASP:OD2	2.47	0.47
1:C:371:VAL:HG22	1:C:393:LEU:HD13	1.97	0.47
1:C:232:ALA:HA	1:C:241:GLU:O	2.15	0.46
1:A:436:ASN:HD22	1:A:449:GLU:HB2	1.80	0.46
1:C:622:ILE:HD11	1:C:631:VAL:HG11	1.97	0.46
1:A:481:MET:HG2	1:A:516:ARG:HG3	1.98	0.45
1:A:549:ALA:HA	1:A:586:THR:HG22	1.97	0.45
2:B:95:THR:HG21	2:B:98:TYR:CZ	2.51	0.45
1:A:430:HIS:CG	1:A:545:LEU:HD13	2.51	0.45
1:C:688:ILE:HG13	1:C:689:VAL:HG23	1.99	0.45
3:C:801:C8E:H82	3:C:801:C8E:H51	1.84	0.45
1:A:583:GLU:HG3	1:A:585:ARG:HG3	2.00	0.44
1:C:628:LEU:HD22	1:C:658:GLU:HB3	1.99	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:536:LEU:HD11	1:C:605:ALA:HB3	2.00	0.43
2:D:86:THR:HG23	2:D:95:THR:HG23	2.00	0.42
1:A:241:GLU:HB3	1:A:261:HIS:HE1	1.84	0.42
1:A:504:LEU:HG	1:A:558:LEU:HD11	2.01	0.42
1:A:228:LEU:HB2	1:A:245:PRO:HB2	2.00	0.42
1:C:418:ASN:OD1	1:C:419:SER:N	2.53	0.42
1:C:554:VAL:HG13	1:C:579:TYR:HB3	2.00	0.42
1:A:240:VAL:HG22	1:A:264:ASN:HB2	2.02	0.42
1:C:486:VAL:O	1:C:509:GLN:HA	2.20	0.41
1:C:719:CYS:SG	1:C:720:CYS:N	2.88	0.41
2:B:29:VAL:HA	2:B:30:PRO:HD3	1.96	0.41
1:C:521:ILE:O	1:C:547:ARG:NH2	2.53	0.41
1:A:583:GLU:HB3	1:A:598:THR:HB	2.00	0.41
1:C:564:ASP:HB3	1:C:570:ARG:HD2	2.02	0.41
2:D:143:ILE:O	2:D:147:MET:HG3	2.20	0.41
1:C:244:LEU:O	1:C:260:PRO:HD2	2.21	0.41
1:C:667:TYR:O	1:C:673:TYR:HB2	2.20	0.41
2:D:36:MET:O	2:D:62:LEU:HA	2.21	0.41
1:A:502:GLN:HA	1:A:562:VAL:HA	2.02	0.40
1:C:424:PRO:HB3	1:C:462:LEU:HD23	2.03	0.40
2:D:64:ASP:OD1	2:D:65:ALA:N	2.50	0.40
1:C:774:ILE:HD12	1:C:774:ILE:HA	1.94	0.40
1:A:694:VAL:HG22	1:A:714:VAL:HG12	2.03	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	548/601 (91%)	510 (93%)	35 (6%)	3 (0%)	34 70
1	C	539/601 (90%)	497 (92%)	38 (7%)	4 (1%)	26 62

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	142/182 (78%)	139 (98%)	3 (2%)	0	100	100
2	D	147/182 (81%)	141 (96%)	6 (4%)	0	100	100
All	All	1376/1566 (88%)	1287 (94%)	82 (6%)	7 (0%)	34	70

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	666	SER
1	A	365	GLN
1	C	629	ASP
1	C	365	GLN
1	C	239	GLY
1	A	221	ASP
1	C	552	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	480/526 (91%)	474 (99%)	6 (1%)	76	93
1	C	473/526 (90%)	464 (98%)	9 (2%)	65	88
2	B	127/156 (81%)	127 (100%)	0	100	100
2	D	131/156 (84%)	130 (99%)	1 (1%)	86	96
All	All	1211/1364 (89%)	1195 (99%)	16 (1%)	76	93

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	406	ASP
1	A	437	LEU
1	A	552	ASN
1	A	589	ASP
1	A	673	TYR

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Mol	Chain	Res	Type
1	A	721	TYR
1	C	290	TYR
1	C	394	ASP
1	C	421	ASN
1	C	425	GLU
1	C	516	ARG
1	C	589	ASP
1	C	677	ILE
1	C	692	TRP
1	C	749	PHE
2	D	28	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
2	B	91	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

8 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	801	-	20,20,20	0.39	0	19,19,19	0.33	0
3	C8E	A	802	-	20,20,20	0.38	0	19,19,19	0.41	0
3	C8E	A	803	-	20,20,20	0.39	0	19,19,19	0.38	0
3	C8E	C	801	-	20,20,20	0.39	0	19,19,19	0.36	0
3	C8E	C	802	-	20,20,20	0.38	0	19,19,19	0.41	0
3	C8E	C	803	-	20,20,20	0.38	0	19,19,19	0.41	0
3	C8E	C	804	-	20,20,20	0.39	0	19,19,19	0.39	0
3	C8E	D	201	-	20,20,20	0.40	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	801	-	-	0/18/18/18	0/0/0/0
3	C8E	A	802	-	-	0/18/18/18	0/0/0/0
3	C8E	A	803	-	-	0/18/18/18	0/0/0/0
3	C8E	C	801	-	-	0/18/18/18	0/0/0/0
3	C8E	C	802	-	-	0/18/18/18	0/0/0/0
3	C8E	C	803	-	-	0/18/18/18	0/0/0/0
3	C8E	C	804	-	-	0/18/18/18	0/0/0/0
3	C8E	D	201	-	-	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.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	801	C8E	2	0
3	C	801	C8E	1	0

5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	552/601 (91%)	0.43	41 (7%) 17 13	14, 62, 120, 154	0
1	C	543/601 (90%)	0.85	100 (18%) 2 1	21, 72, 157, 211	0
2	B	146/182 (80%)	0.54	13 (8%) 12 8	21, 63, 126, 164	0
2	D	149/182 (81%)	0.41	9 (6%) 25 20	29, 62, 134, 180	0
All	All	1390/1566 (88%)	0.60	163 (11%) 6 4	14, 65, 144, 211	0

All (163) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	753	ILE	8.0
1	C	688	ILE	7.3
1	C	723	ILE	6.3
1	C	301	HIS	6.2
1	A	722	ALA	6.1
1	C	267	GLY	6.0
1	C	694	VAL	5.8
1	C	275	PHE	5.5
1	C	269	VAL	5.5
2	B	29	VAL	5.5
1	C	230	PRO	5.2
1	C	716	TYR	5.2
1	C	304	ASP	5.2
1	C	717	ASN	4.9
1	C	751	ILE	4.9
1	C	689	VAL	4.9
1	C	244	LEU	4.7
1	C	344	SER	4.7
2	B	69	ARG	4.7
1	C	382	GLU	4.7
1	A	720	CYS	4.6

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Mol	Chain	Res	Type	RSRZ
1	C	695	VAL	4.6
1	C	312	PHE	4.6
2	D	27	THR	4.5
1	C	306	ASN	4.5
1	C	692	TRP	4.5
1	C	693	SER	4.4
1	C	229	ILE	4.4
2	B	22	HIS	4.3
2	B	20	GLY	4.3
1	C	690	ASP	4.3
1	A	596	ASN	4.2
2	B	23	LEU	4.2
1	C	240	VAL	4.1
2	D	26	THR	4.1
1	C	228	LEU	4.1
2	B	68	LEU	4.1
1	C	263	MET	4.1
1	C	719	CYS	4.0
1	C	262	TYR	4.0
1	C	255	ASP	4.0
1	C	725	LEU	4.0
1	C	302	SER	4.0
1	C	232	ALA	4.0
1	A	723	ILE	3.9
1	C	256	ALA	3.9
1	C	721	TYR	3.9
1	C	686	TRP	3.9
1	C	718	SER	3.8
1	C	720	CYS	3.7
1	C	752	ASN	3.7
1	C	271	TRP	3.7
1	A	690	ASP	3.6
2	D	29	VAL	3.6
1	C	288	PHE	3.6
1	C	242	PHE	3.5
1	C	714	VAL	3.5
1	C	724	ARG	3.5
1	C	687	PRO	3.5
1	C	781	LEU	3.4
2	B	21	TRP	3.4
1	A	492	PHE	3.4
1	C	260	PRO	3.4

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Mol	Chain	Res	Type	RSRZ
1	C	249	ASN	3.4
1	C	264	ASN	3.3
1	C	712	LEU	3.3
1	C	297	TYR	3.3
1	A	565	ALA	3.3
2	B	33	MET	3.3
2	D	33	MET	3.3
1	A	220	GLY	3.3
1	C	276	ARG	3.2
1	C	749	PHE	3.2
1	C	293	SER	3.2
1	C	684	ALA	3.2
1	C	247	TYR	3.2
2	D	68	LEU	3.1
1	C	227	PHE	3.1
1	C	722	ALA	3.1
1	A	595	ASN	3.1
2	D	67	THR	3.1
2	B	70	LYS	3.1
1	C	231	ASN	3.1
1	A	472	ASP	3.0
1	A	691	ARG	3.0
1	A	714	VAL	3.0
1	A	688	ILE	3.0
2	B	24	ARG	3.0
1	C	467	ALA	3.0
1	C	736	SER	3.0
1	C	305	SER	3.0
1	C	342	PHE	2.9
1	C	729	ARG	2.9
1	C	333	VAL	2.9
1	C	691	ARG	2.9
1	A	719	CYS	2.9
1	A	615	GLU	2.9
1	C	682	MET	2.9
1	C	246	TYR	2.9
1	C	727	TYR	2.8
1	C	233	LYS	2.8
1	C	258	ILE	2.8
2	B	71	ASP	2.8
1	C	252	PRO	2.8
1	A	712	LEU	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	303	SER	2.7
1	C	298	GLU	2.7
1	C	234	TYR	2.7
1	A	305	SER	2.7
1	C	698	TYR	2.7
1	C	286	THR	2.6
1	C	346	TYR	2.6
1	C	331	THR	2.6
1	C	685	SER	2.6
2	D	116	TYR	2.6
1	C	253	ASN	2.6
1	C	612	ILE	2.6
1	A	739	ASN	2.6
1	A	225	SER	2.5
1	C	381	ARG	2.5
1	A	234	TYR	2.5
1	C	250	ILE	2.5
1	C	311	LEU	2.5
1	C	248	TRP	2.5
1	C	257	THR	2.5
1	C	254	PHE	2.4
1	C	746	ASP	2.4
1	A	702	THR	2.4
2	B	31	THR	2.4
1	C	700	TYR	2.4
1	A	721	TYR	2.4
1	C	259	THR	2.4
1	A	737	ASN	2.3
1	A	717	ASN	2.3
1	C	265	LYS	2.3
1	A	740	GLY	2.3
2	D	28	LYS	2.3
1	A	265	LYS	2.3
1	A	753	ILE	2.3
1	A	224	ARG	2.2
1	A	594	GLU	2.2
1	A	221	ASP	2.2
1	A	678	SER	2.2
1	A	495	ASP	2.2
1	C	310	TRP	2.2
1	C	251	ALA	2.2
1	A	707	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	C	261	HIS	2.1
1	A	724	ARG	2.1
1	C	498	GLU	2.1
1	A	708	ALA	2.1
2	B	108	VAL	2.1
1	C	291	LEU	2.1
1	A	686	TRP	2.1
2	D	111	PRO	2.1
1	A	500	PHE	2.1
1	C	266	ARG	2.1
1	A	689	VAL	2.1
1	C	769	MET	2.1
1	A	602	LEU	2.0
1	A	223	ARG	2.0
1	C	697	ALA	2.0
1	A	752	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
3	C8E	D	201	21/21	0.88	0.38	9.42	19,62,96,105	0
3	C8E	C	803	21/21	0.86	0.31	5.65	42,80,130,132	0
3	C8E	C	804	21/21	0.92	0.34	5.25	28,45,61,65	0
3	C8E	A	802	21/21	0.89	0.39	4.32	22,53,94,110	0
3	C8E	A	801	21/21	0.88	0.30	3.41	31,69,97,99	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	C8E	C	802	21/21	0.92	0.26	3.13	22,48,72,79	0
3	C8E	C	801	21/21	0.93	0.27	2.29	16,46,75,84	0
3	C8E	A	803	21/21	0.76	0.29	1.83	32,74,93,100	0

6.5 Other polymers [i](#)

There are no such residues in this entry.