



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 11:29 PM GMT

PDB ID : 4HMK
Title : Crystal structure of LeuT-E290S with bound Br
Authors : Kantcheva, A.K.; Quick, M.; Shi, L.; Winther, A.M.L.; Stolzenberg, S.; Weinstein, H.; Javitch, J.A.; Nissen, P.
Deposited on : 2012-10-18
Resolution : 3.00 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

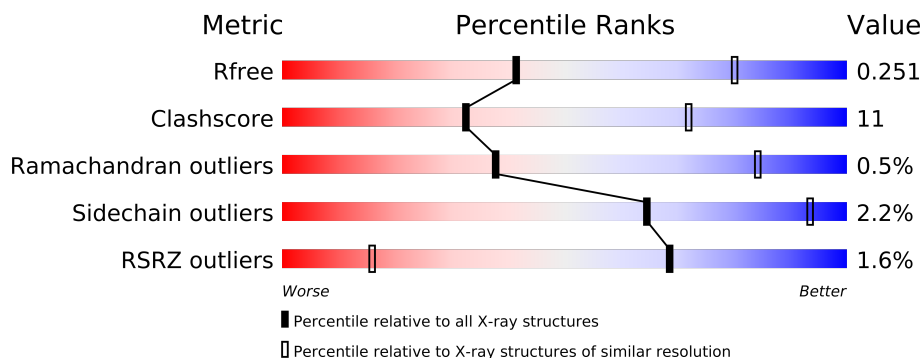
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 3.00 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	66092	1216 (3.00-3.00)
Clashscore	79885	1594 (3.00-3.00)
Ramachandran outliers	78287	1537 (3.00-3.00)
Sidechain outliers	78261	1540 (3.00-3.00)
RSRZ outliers	66119	1217 (3.00-3.00)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	513	
1	B	513	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
3	NA	A	602	-	X
3	NA	B	604	-	X
4	BR	A	613	-	X
5	BOG	A	606	-	X
5	BOG	A	608	-	X
5	BOG	A	609	-	X

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Mol	Type	Chain	Res	Geometry	Electron density
5	BOG	A	610	-	X
5	BOG	B	602	-	X
5	BOG	B	608	-	X
5	BOG	B	609	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8240 atoms, of which 0 are hydrogen and 0 are deuterium.

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.

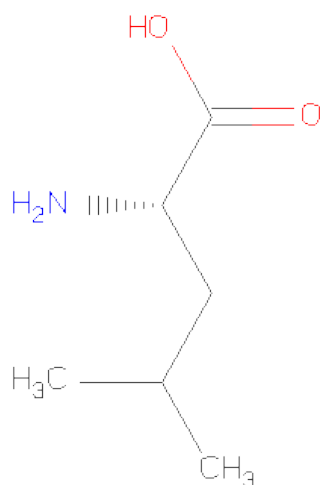
- Molecule 1 is a protein called Transporter.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	501	Total	C	N	O	S	0	0	0
			3987	2700	626	650	11			
1	B	506	Total	C	N	O	S	0	0	0
			4027	2728	632	655	12			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	290	SER	GLU	ENGINEERED MUTATION	UNP O67854
B	290	SER	GLU	ENGINEERED MUTATION	UNP O67854

- Molecule 2 is LEUCINE (three-letter code: LEU) (formula: C₆H₁₃NO₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			9	6	1	2		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	B	1	Total	C	N	O	0	0
			9	6	1	2		

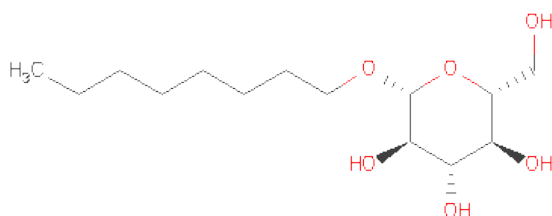
- Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	2	Total	Na	0	0
			2	2		
3	A	2	Total	Na	0	0
			2	2		

- Molecule 4 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	5	Total	Br	0	0
			5	5		
4	A	6	Total	Br	0	0
			6	6		

- Molecule 5 is SUGAR (B-OCTYLGLUCOSIDE) (three-letter code: BOG) (formula: C₁₄H₂₈O₆).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			20	14	6		
5	A	1	Total	C	O	0	0
			20	14	6		

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 20 14 6	0	0
5	A	1	Total C O 20 14 6	0	0
5	A	1	Total C O 20 14 6	0	0
5	B	1	Total C O 20 14 6	0	0
5	B	1	Total C O 20 14 6	0	0
5	B	1	Total C O 20 14 6	0	0
5	B	1	Total C O 20 14 6	0	0

- Molecule 6 is water.

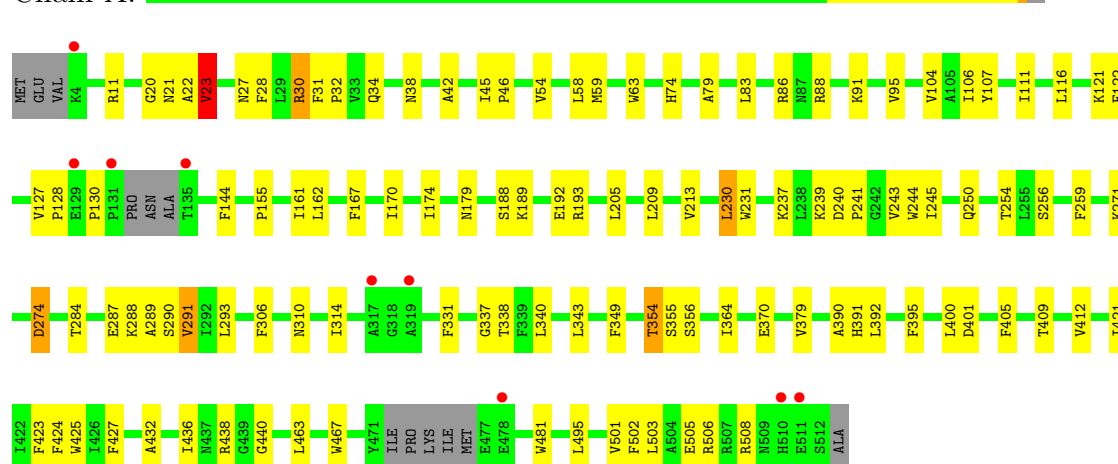
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	7	Total O 7 7	0	0
6	B	6	Total O 6 6	0	0

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.

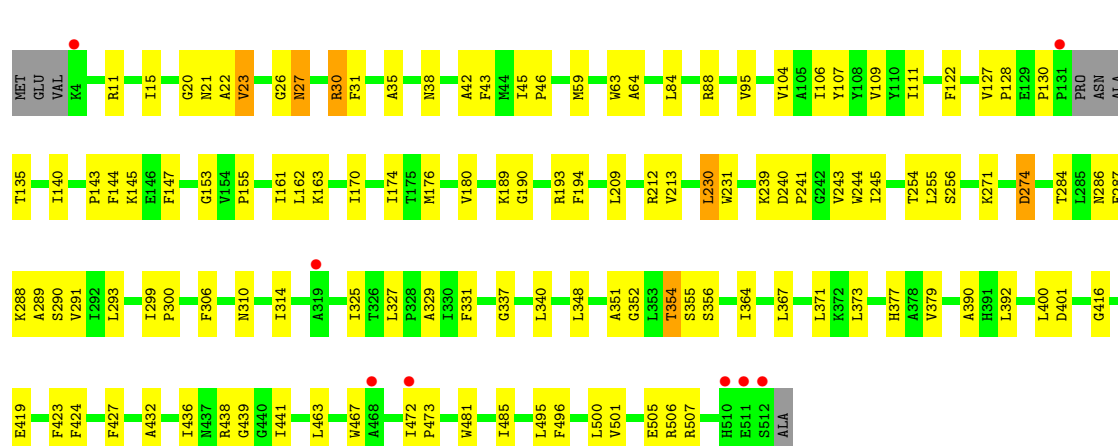
• Molecule 1: Transporter

Chain A:



• Molecule 1: Transporter

Chain B:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	84.14Å 92.32Å 87.21Å 90.00° 94.25° 90.00°	Depositor
Resolution (Å)	63.30 – 3.00 63.30 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.7 (63.30-3.00) 99.7 (63.30-3.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.61 (at 3.01Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.6.4_486)	Depositor
R, R_{free}	0.205 , 0.261 0.201 , 0.251	Depositor DCC
R_{free} test set	1337 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	63.2	Xtriage
Anisotropy	0.264	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 43.0	EDS
Estimated twinning fraction	0.024 for l,-k,h	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 26769 reflections	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	8240	wwPDB-VP
Average B, all atoms (Å ²)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 17.27% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA, BR, BOG

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	0.42	1/4111 (0.0%)	0.56	1/5596 (0.0%)
1	B	0.40	0/4153	0.55	1/5654 (0.0%)
All	All	0.41	1/8264 (0.0%)	0.55	2/11250 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	23	VAL	CB-CG2	-5.10	1.42	1.52

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	351	ALA	C-N-CA	-5.57	110.60	122.30
1	A	354	THR	OG1-CB-CG2	-5.51	97.33	110.00

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3987	0	4071	95	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	4027	0	4123	92	0
2	A	9	0	10	1	0
2	B	9	0	10	2	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	6	0	0	4	0
4	B	5	0	0	7	0
5	A	100	0	139	28	0
5	B	80	0	111	10	0
6	A	7	0	0	0	0
6	B	6	0	0	1	0
All	All	8240	0	8464	189	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (189) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
5:A:609:BOG:H5	1:B:507:ARG:HA	1.47	0.96
1:A:86:ARG:HH12	5:A:610:BOG:H5'2	1.30	0.93
5:A:608:BOG:H8'3	5:B:601:BOG:H8'2	1.53	0.91
1:A:27:ASN:HD22	1:A:290:SER:HB3	1.38	0.88
1:A:11:ARG:HH22	1:A:274:ASP:HB3	1.41	0.86
5:A:608:BOG:H6'2	5:B:601:BOG:H6'1	1.62	0.81
1:B:11:ARG:HH22	1:B:274:ASP:HB3	1.43	0.81
1:B:364:ILE:HD13	1:B:379:VAL:HG22	1.64	0.78
1:A:86:ARG:NH1	5:A:610:BOG:H5'2	1.97	0.78
1:A:161:ILE:HD12	5:B:601:BOG:H5'1	1.68	0.75
5:A:609:BOG:H5'2	1:B:371:LEU:O	1.88	0.74
1:A:364:ILE:HD13	1:A:379:VAL:HG22	1.70	0.73
1:A:392:LEU:HD13	5:A:608:BOG:H8'1	1.72	0.71
1:B:155:PRO:HD3	1:B:162:LEU:HD23	1.74	0.70
1:A:271:LYS:O	1:A:438:ARG:NH1	2.24	0.70
1:B:135:THR:N	6:B:703:HOH:O	2.26	0.69
1:B:27:ASN:HD22	1:B:290:SER:HB3	1.58	0.69
1:A:189:LYS:HA	1:A:193:ARG:HG3	1.77	0.67
1:A:83:LEU:HD21	5:A:610:BOG:H4'2	1.76	0.67
1:B:256:SER:HG	2:B:603:LEU:N	1.95	0.65
5:A:609:BOG:H8'1	4:B:611:BR:BR	2.53	0.64
1:A:155:PRO:HD3	1:A:162:LEU:HD23	1.79	0.64
1:A:245:ILE:HG12	1:A:463:LEU:HD12	1.79	0.64
1:A:256:SER:HG	2:A:601:LEU:N	1.96	0.63

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:367:LEU:HB3	1:B:373:LEU:HD22	1.81	0.63
1:A:11:ARG:HH22	1:A:274:ASP:CB	2.11	0.61
1:B:45:ILE:HB	1:B:46:PRO:HD3	1.82	0.61
1:A:88:ARG:NH1	1:A:91:LYS:HE3	2.15	0.61
1:B:11:ARG:HH22	1:B:274:ASP:CB	2.15	0.60
1:B:11:ARG:NH2	1:B:274:ASP:HB3	2.15	0.59
1:B:189:LYS:HA	1:B:193:ARG:HG3	1.83	0.59
1:B:392:LEU:HD13	5:B:609:BOG:H8'1	1.84	0.58
1:B:241:PRO:O	1:B:245:ILE:HG13	2.03	0.58
1:A:503:LEU:HD22	5:A:609:BOG:H4'2	1.85	0.58
1:A:338:THR:HB	4:A:613:BR:BR	2.58	0.58
1:A:23:VAL:HG12	1:A:23:VAL:O	2.04	0.57
1:A:111:ILE:HD11	1:A:400:LEU:HD11	1.86	0.57
1:A:45:ILE:HB	1:A:46:PRO:HD3	1.86	0.57
1:B:31:PHE:CD1	4:B:606:BR:BR	3.13	0.57
1:A:11:ARG:NH2	1:A:274:ASP:HB3	2.15	0.56
1:B:245:ILE:HG12	1:B:463:LEU:HD12	1.88	0.56
1:A:331:PHE:O	1:A:337:GLY:HA3	2.06	0.56
1:A:88:ARG:HA	1:A:88:ARG:HH11	1.69	0.56
1:A:31:PHE:HB2	4:A:604:BR:BR	2.61	0.55
1:B:377:HIS:ND1	4:B:611:BR:BR	2.87	0.55
1:B:31:PHE:HB2	4:B:606:BR:BR	2.62	0.54
1:A:395:PHE:HB3	5:B:601:BOG:H5'2	1.90	0.54
1:B:23:VAL:HG12	1:B:23:VAL:O	2.06	0.54
1:A:34:GLN:NE2	4:A:611:BR:BR	2.95	0.54
1:A:506:ARG:HG3	5:A:609:BOG:H8'3	1.89	0.54
1:B:212:ARG:HG3	1:B:212:ARG:HH11	1.73	0.54
1:A:74:HIS:NE2	5:A:610:BOG:H3'1	2.24	0.53
1:A:130:PRO:HG3	1:A:144:PHE:HE2	1.72	0.53
1:B:254:THR:OG1	4:B:606:BR:BR	2.79	0.53
1:A:27:ASN:ND2	1:A:290:SER:HB3	2.18	0.52
1:A:241:PRO:O	1:A:245:ILE:HG13	2.09	0.52
1:B:289:ALA:O	1:B:293:LEU:HB2	2.10	0.52
1:A:27:ASN:HB3	1:A:290:SER:HB3	1.91	0.51
5:A:609:BOG:H8'2	1:B:373:LEU:HD12	1.93	0.51
1:A:167:PHE:HE1	5:A:606:BOG:H6'2	1.76	0.50
1:B:255:LEU:HD23	1:B:286:ASN:HD22	1.75	0.50
1:A:30:ARG:HD3	4:A:611:BR:BR	2.67	0.50
1:A:239:LYS:HD2	1:A:239:LYS:N	2.25	0.50
1:B:95:VAL:HG22	1:B:501:VAL:HG21	1.93	0.50
1:B:271:LYS:O	1:B:438:ARG:NH1	2.41	0.50
1:A:370:GLU:OE1	5:A:610:BOG:H61	2.11	0.50

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:46:PRO:HB3	1:A:244:TRP:CE3	2.47	0.50
1:A:88:ARG:HD3	1:A:91:LYS:HD2	1.94	0.49
1:A:167:PHE:HD1	5:A:606:BOG:H4'2	1.77	0.49
1:B:239:LYS:HD2	1:B:239:LYS:N	2.27	0.49
1:B:88:ARG:HH11	1:B:88:ARG:HA	1.77	0.49
1:B:284:THR:HG22	1:B:288:LYS:HE3	1.94	0.49
1:B:230:LEU:HB3	1:B:231:TRP:CD1	2.47	0.49
1:A:506:ARG:CB	5:A:609:BOG:H6'2	2.43	0.49
1:A:423:PHE:O	1:A:427:PHE:HB3	2.12	0.49
1:A:495:LEU:CB	5:B:602:BOG:H61	2.42	0.49
1:B:432:ALA:O	1:B:436:ILE:HG13	2.13	0.48
1:A:495:LEU:HB2	5:B:602:BOG:H61	1.96	0.48
1:B:240:ASP:HB3	1:B:243:VAL:HG23	1.96	0.48
1:B:27:ASN:HB3	1:B:290:SER:HB3	1.96	0.48
1:B:287:GLU:O	1:B:291:VAL:HG13	2.13	0.48
1:B:416:GLY:O	1:B:419:GLU:HG2	2.14	0.48
1:B:127:VAL:HB	1:B:128:PRO:HD2	1.96	0.48
1:A:79:ALA:HA	5:A:610:BOG:H2	1.96	0.47
1:B:46:PRO:HB3	1:B:244:TRP:CE3	2.49	0.47
1:B:104:VAL:HA	1:B:107:TYR:CE2	2.49	0.47
1:B:505:GLU:HG3	1:B:506:ARG:NH1	2.30	0.47
1:B:495:LEU:HB3	5:B:602:BOG:H8'2	1.95	0.47
1:A:179:ASN:OD1	1:A:356:SER:OG	2.30	0.47
1:B:30:ARG:HD3	4:B:610:BR:BR	2.70	0.47
1:B:245:ILE:HG12	1:B:463:LEU:CD1	2.45	0.47
1:A:130:PRO:HG3	1:A:144:PHE:CE2	2.49	0.47
1:B:163:LYS:NZ	5:B:609:BOG:H61	2.30	0.47
1:A:505:GLU:HG3	1:A:506:ARG:NH1	2.30	0.47
1:A:240:ASP:HB3	1:A:243:VAL:HG23	1.96	0.46
1:B:231:TRP:HH2	1:B:291:VAL:HB	1.80	0.46
1:A:250:GLN:O	1:A:254:THR:HG22	2.15	0.46
1:B:11:ARG:O	1:B:15:ILE:HG12	2.16	0.46
1:A:506:ARG:HB2	5:A:609:BOG:H6'2	1.97	0.46
1:A:122:PHE:CD1	1:A:127:VAL:HG11	2.50	0.46
1:B:122:PHE:CD1	1:B:127:VAL:HG11	2.51	0.46
5:A:609:BOG:C8'	4:B:611:BR:BR	3.18	0.46
1:A:306:PHE:HB3	1:A:310:ASN:HB2	1.98	0.46
1:B:111:ILE:HD11	1:B:400:LEU:HD11	1.98	0.46
1:A:245:ILE:HG12	1:A:463:LEU:CD1	2.46	0.45
1:A:20:GLY:O	1:A:21:ASN:C	2.51	0.45
1:A:287:GLU:O	1:A:291:VAL:HG13	2.17	0.45
1:A:59:MET:O	1:A:63:TRP:HD1	1.99	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:331:PHE:O	1:B:337:GLY:HA3	2.16	0.45
1:A:432:ALA:O	1:A:436:ILE:HG13	2.16	0.45
1:A:205:LEU:HD22	1:A:343:LEU:HD22	1.98	0.45
5:A:609:BOG:H5	1:B:507:ARG:CA	2.34	0.45
1:A:83:LEU:CD2	5:A:610:BOG:H4'2	2.44	0.45
1:B:153:GLY:HA3	1:B:163:LYS:O	2.17	0.45
1:B:255:LEU:HD23	1:B:286:ASN:ND2	2.31	0.45
1:B:306:PHE:HB3	1:B:310:ASN:HB2	1.98	0.45
1:B:472:ILE:N	1:B:473:PRO:HD2	2.32	0.45
1:B:140:ILE:O	1:B:143:PRO:HD2	2.17	0.45
1:A:128:PRO:HG2	1:A:144:PHE:CE1	2.50	0.45
1:B:255:LEU:CD2	1:B:286:ASN:HD22	2.29	0.45
1:B:109:VAL:HG22	1:B:356:SER:HB2	1.99	0.44
1:A:161:ILE:HD11	1:B:481:TRP:CE3	2.51	0.44
1:A:106:ILE:HG23	1:A:390:ALA:HA	1.99	0.44
1:B:310:ASN:O	1:B:314:ILE:HG13	2.17	0.44
1:A:340:LEU:HA	1:A:340:LEU:HD12	1.75	0.44
1:A:502:PHE:O	1:A:506:ARG:HG2	2.17	0.44
1:A:230:LEU:HB3	1:A:231:TRP:CD1	2.53	0.44
1:A:259:PHE:CZ	1:A:412:VAL:HG11	2.53	0.44
1:A:237:LYS:HB3	1:A:243:VAL:HG21	2.00	0.44
1:A:167:PHE:CE1	5:A:606:BOG:H6'2	2.53	0.44
1:A:104:VAL:HA	1:A:107:TYR:CE2	2.53	0.44
1:B:423:PHE:O	1:B:427:PHE:HB3	2.17	0.44
1:B:38:ASN:O	1:B:42:ALA:HB3	2.17	0.44
1:B:35:ALA:HA	1:B:43:PHE:CD2	2.53	0.43
1:B:46:PRO:HG2	1:B:243:VAL:HG12	1.99	0.43
1:B:170:ILE:O	1:B:174:ILE:HG13	2.18	0.43
1:B:144:PHE:CE2	1:B:329:ALA:HA	2.53	0.43
1:A:370:GLU:HB3	5:A:610:BOG:H62	2.01	0.43
1:A:274:ASP:OD1	1:A:440:GLY:HA2	2.19	0.43
1:B:26:GLY:HA3	2:B:603:LEU:OXT	2.18	0.43
1:B:145:LYS:HA	1:B:325:ILE:HD13	2.01	0.43
1:B:106:ILE:HG23	1:B:390:ALA:HA	2.00	0.43
1:A:161:ILE:HD11	1:B:481:TRP:CZ3	2.54	0.43
1:A:505:GLU:HB2	1:A:508:ARG:HH21	1.84	0.43
1:A:245:ILE:HG23	1:A:463:LEU:HD11	2.00	0.43
1:B:153:GLY:HA3	1:B:163:LYS:C	2.39	0.43
1:B:340:LEU:HD12	1:B:340:LEU:HA	1.83	0.42
1:B:348:LEU:O	1:B:352:GLY:N	2.52	0.42
1:A:289:ALA:O	1:A:293:LEU:HB2	2.19	0.42
1:B:194:PHE:HE2	1:B:354:THR:HG23	1.85	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:128:PRO:HD3	1:B:147:PHE:CD1	2.54	0.42
1:A:481:TRP:CE3	1:B:161:ILE:HD11	2.55	0.42
1:A:391:HIS:HD2	5:A:607:BOG:H3'1	1.85	0.42
1:B:11:ARG:HH21	1:B:441:ILE:HD12	1.85	0.42
1:B:84:LEU:HB3	1:B:427:PHE:HB2	2.02	0.42
1:A:481:TRP:CZ3	1:B:161:ILE:HD11	2.55	0.42
1:B:496:PHE:CZ	1:B:500:LEU:HD11	2.55	0.42
1:A:188:SER:O	1:A:192:GLU:HB3	2.20	0.42
5:A:609:BOG:H1'1	1:B:507:ARG:HB2	2.01	0.41
1:B:299:ILE:HB	1:B:300:PRO:HD3	2.02	0.41
1:B:194:PHE:CE2	1:B:354:THR:HG23	2.55	0.41
5:A:607:BOG:H6'1	5:A:607:BOG:H3'2	1.90	0.41
1:A:405:PHE:O	1:A:409:THR:HB	2.19	0.41
1:A:54:VAL:O	1:A:58:LEU:HB2	2.20	0.41
5:A:610:BOG:H3'2	5:A:610:BOG:H7'2	2.01	0.41
1:A:284:THR:HG22	1:A:288:LYS:HE3	2.02	0.41
1:A:116:LEU:HB2	1:A:349:PHE:CD1	2.56	0.41
1:A:310:ASN:O	1:A:314:ILE:HG13	2.19	0.41
1:B:176:MET:O	1:B:180:VAL:HG23	2.20	0.41
1:B:485:ILE:HG12	5:B:601:BOG:H8'3	2.03	0.41
1:A:46:PRO:HG2	1:A:243:VAL:HG12	2.02	0.41
1:B:64:ALA:HB2	1:B:436:ILE:HG12	2.03	0.41
1:B:209:LEU:O	1:B:213:VAL:HG23	2.21	0.41
1:A:421:ILE:O	1:A:425:TRP:HB2	2.21	0.41
1:B:274:ASP:HB2	1:B:439:GLY:O	2.21	0.41
1:A:95:VAL:HG22	1:A:501:VAL:HG21	2.03	0.41
1:B:327:LEU:HA	1:B:327:LEU:HD12	1.83	0.41
1:B:212:ARG:HG3	1:B:212:ARG:NH1	2.36	0.41
1:B:20:GLY:O	1:B:21:ASN:C	2.55	0.40
1:A:161:ILE:HA	1:A:161:ILE:HD13	1.87	0.40
1:A:28:PHE:O	1:A:32:PRO:HG2	2.21	0.40
1:A:170:ILE:O	1:A:174:ILE:HG13	2.21	0.40
1:A:38:ASN:O	1:A:42:ALA:HB3	2.21	0.40
1:B:59:MET:O	1:B:63:TRP:HD1	2.04	0.40
1:A:23:VAL:O	1:A:23:VAL:CG1	2.67	0.40
1:A:205:LEU:CD2	1:A:343:LEU:HD22	2.51	0.40
1:B:130:PRO:HG3	1:B:144:PHE:HE2	1.87	0.40
1:A:209:LEU:O	1:A:213:VAL:HG23	2.21	0.40
1:A:121:LYS:HD2	1:A:167:PHE:CE2	2.55	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution. The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	495/513 (96%)	475 (96%)	18 (4%)	2 (0%)	43	87
1	B	502/513 (98%)	469 (93%)	30 (6%)	3 (1%)	33	81
All	All	997/1026 (97%)	944 (95%)	48 (5%)	5 (0%)	38	84

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	23	VAL
1	A	23	VAL
1	B	22	ALA
1	A	22	ALA
1	B	190	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	405/415 (98%)	396 (98%)	9 (2%)	64	93
1	B	410/415 (99%)	401 (98%)	9 (2%)	64	93
All	All	815/830 (98%)	797 (98%)	18 (2%)	64	93

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	30	ARG
1	A	230	LEU
1	A	274	ASP
1	A	291	VAL

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Mol	Chain	Res	Type
1	A	354	THR
1	A	355	SER
1	A	401	ASP
1	A	424	PHE
1	A	467	TRP
1	B	27	ASN
1	B	30	ARG
1	B	230	LEU
1	B	274	ASP
1	B	354	THR
1	B	355	SER
1	B	401	ASP
1	B	424	PHE
1	B	467	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

Of 26 ligands modelled in this entry, 15 are monoatomic - leaving 11 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	LEU	A	601	3	8,8,8	1.05	1 (12%)	10,10,10	1.82	3 (30%)
5	BOG	A	606	-	20,20,20	1.30	3 (15%)	25,25,25	1.22	3 (12%)
5	BOG	A	607	-	20,20,20	1.28	3 (15%)	25,25,25	1.44	3 (12%)
5	BOG	A	608	-	20,20,20	1.28	3 (15%)	25,25,25	1.24	3 (12%)
5	BOG	A	609	-	20,20,20	1.40	3 (15%)	25,25,25	1.26	4 (16%)
5	BOG	A	610	-	20,20,20	2.32	7 (35%)	25,25,25	2.21	6 (24%)
5	BOG	B	601	-	20,20,20	1.32	3 (15%)	25,25,25	1.31	3 (12%)
5	BOG	B	602	-	20,20,20	1.72	5 (25%)	25,25,25	2.73	9 (36%)
2	LEU	B	603	3	8,8,8	0.89	1 (12%)	10,10,10	1.98	3 (30%)
5	BOG	B	608	-	20,20,20	1.37	3 (15%)	25,25,25	1.35	2 (8%)
5	BOG	B	609	-	20,20,20	1.34	3 (15%)	25,25,25	1.04	2 (8%)

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	LEU	A	601	3	-	0/8/8/8	0/0/0/0
5	BOG	A	606	-	-	0/11/31/31	0/1/1/1
5	BOG	A	607	-	-	0/11/31/31	0/1/1/1
5	BOG	A	608	-	-	0/11/31/31	0/1/1/1
5	BOG	A	609	-	-	0/11/31/31	0/1/1/1
5	BOG	A	610	-	-	0/11/31/31	0/1/1/1
5	BOG	B	601	-	-	0/11/31/31	0/1/1/1
5	BOG	B	602	-	-	0/11/31/31	0/1/1/1
2	LEU	B	603	3	-	0/8/8/8	0/0/0/0
5	BOG	B	608	-	-	0/11/31/31	0/1/1/1
5	BOG	B	609	-	-	0/11/31/31	0/1/1/1

All (35) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	A	610	BOG	O5-C1	5.04	1.54	1.41
5	A	610	BOG	O1-C1	-4.61	1.31	1.40
5	A	610	BOG	O2-C2	-4.59	1.31	1.43
5	B	602	BOG	O5-C1	4.50	1.53	1.41
5	B	601	BOG	O5-C1	4.11	1.52	1.41
5	B	608	BOG	O5-C1	4.04	1.52	1.41
5	A	609	BOG	O5-C1	3.93	1.51	1.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	B	609	BOG	O5-C1	3.92	1.51	1.41
5	A	606	BOG	O5-C1	3.87	1.51	1.41
5	A	607	BOG	O5-C1	3.87	1.51	1.41
5	A	610	BOG	C3-C2	-3.78	1.42	1.52
5	A	608	BOG	O5-C1	3.56	1.51	1.41
5	B	602	BOG	C3-C2	-3.13	1.44	1.52
5	A	609	BOG	O1-C1	-2.92	1.34	1.40
5	B	602	BOG	O3-C3	-2.82	1.36	1.43
2	A	601	LEU	OXT-C	-2.81	1.20	1.30
5	A	608	BOG	O1-C1	-2.68	1.35	1.40
5	A	610	BOG	C4-C3	-2.67	1.45	1.52
5	B	609	BOG	O1-C1	-2.67	1.35	1.40
5	A	610	BOG	C1-C2	-2.57	1.44	1.52
5	A	608	BOG	C3-C2	-2.51	1.45	1.52
5	A	606	BOG	O1-C1	-2.49	1.35	1.40
5	A	609	BOG	C3-C2	-2.43	1.45	1.52
2	B	603	LEU	OXT-C	-2.43	1.21	1.30
5	B	608	BOG	C3-C2	-2.42	1.46	1.52
5	B	609	BOG	C3-C2	-2.40	1.46	1.52
5	B	602	BOG	C4-C5	-2.35	1.47	1.53
5	A	606	BOG	C3-C2	-2.35	1.46	1.52
5	B	608	BOG	O1-C1	-2.31	1.35	1.40
5	B	602	BOG	C4-C3	-2.29	1.46	1.52
5	A	610	BOG	O5-C5	-2.29	1.38	1.44
5	A	607	BOG	O1-C1	-2.21	1.36	1.40
5	A	607	BOG	C3-C2	-2.19	1.46	1.52
5	B	601	BOG	O1-C1	-2.19	1.36	1.40
5	B	601	BOG	C3-C2	-2.17	1.46	1.52

All (41) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	610	BOG	O2-C2-C1	-7.50	93.69	110.04
5	B	602	BOG	C1'-O1-C1	7.12	126.76	113.96
5	B	602	BOG	O3-C3-C4	-5.33	98.41	110.35
5	B	602	BOG	O3-C3-C2	-5.28	98.52	110.35
5	A	607	BOG	C1'-O1-C1	4.76	122.52	113.96
5	B	602	BOG	C6-C5-C4	-4.72	101.60	113.00
5	B	608	BOG	C1'-O1-C1	4.70	122.42	113.96
2	B	603	LEU	OXT-C-O	-4.07	114.87	124.07
2	A	601	LEU	OXT-C-O	-4.05	114.91	124.07
5	B	602	BOG	O5-C5-C6	3.83	115.74	106.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	610	BOG	C3-C4-C5	-3.70	103.59	110.20
5	A	610	BOG	O2-C2-C3	-3.69	102.08	110.35
5	A	608	BOG	O1-C1-C2	3.60	112.76	108.18
5	A	610	BOG	C1-C2-C3	3.27	116.36	110.00
2	B	603	LEU	C-CA-N	3.27	114.77	109.36
5	A	606	BOG	O1-C1-C2	3.24	112.30	108.18
5	A	609	BOG	O5-C5-C6	3.24	114.29	106.34
5	B	601	BOG	C1'-O1-C1	2.96	119.29	113.96
5	B	609	BOG	C1'-O1-C1	2.92	119.21	113.96
5	A	607	BOG	O1-C1-C2	2.80	111.74	108.18
5	B	602	BOG	O1-C1-C2	2.77	111.70	108.18
5	A	608	BOG	C1'-O1-C1	2.71	118.84	113.96
5	A	609	BOG	C1'-O1-C1	2.69	118.80	113.96
5	A	609	BOG	C1-O5-C5	-2.68	108.52	113.73
5	B	602	BOG	O2-C2-C1	2.62	115.75	110.04
5	B	602	BOG	O6-C6-C5	2.62	120.37	111.36
5	A	610	BOG	O4-C4-C5	2.43	115.69	109.28
2	A	601	LEU	CB-CA-C	-2.28	104.51	111.08
5	A	609	BOG	O5-C5-C4	2.25	113.93	109.76
5	B	601	BOG	O1-C1'-C2'	2.25	118.66	109.87
5	B	602	BOG	O1-C1'-C2'	2.23	118.58	109.87
5	B	608	BOG	C6-C5-C4	-2.18	107.73	113.00
5	B	601	BOG	C4-C3-C2	-2.16	106.81	110.82
5	A	608	BOG	O6-C6-C5	2.16	118.80	111.36
5	A	607	BOG	C1-O5-C5	-2.13	109.58	113.73
2	B	603	LEU	CB-CA-C	-2.13	104.96	111.08
5	A	610	BOG	O1-C1'-C2'	2.08	117.99	109.87
5	A	606	BOG	C1'-O1-C1	2.08	117.69	113.96
5	A	606	BOG	O1-C1'-C2'	2.05	117.86	109.87
5	B	609	BOG	O6-C6-C5	2.04	118.39	111.36
2	A	601	LEU	C-CA-N	2.02	112.71	109.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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	501/513 (97%)	-0.06	9 (1%)	65 14	34, 50, 87, 110	0
1	B	506/513 (98%)	-0.10	8 (1%)	68 16	34, 52, 90, 110	0
All	All	1007/1026 (98%)	-0.08	17 (1%)	68 15	34, 51, 88, 110	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	131	PRO	5.9
1	B	511	GLU	4.2
1	A	511	GLU	2.8
1	A	478	GLU	2.7
1	B	468	ALA	2.7
1	B	510	HIS	2.6
1	B	512	SER	2.4
1	B	4	LYS	2.4
1	A	129	GLU	2.2
1	A	317	ALA	2.1
1	B	472	ILE	2.1
1	A	510	HIS	2.1
1	B	319	ALA	2.1
1	A	4	LYS	2.1
1	A	319	ALA	2.1
1	A	135	THR	2.0
1	B	131	PRO	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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	RSR	LLDF	B-factors(Å ²)	Q<0.9
5	BOG	B	609	20/20	0.50	18.68	34,71,80,81	20
4	BR	A	613	1/1	0.55	17.46	237,237,237,237	0
5	BOG	B	602	20/20	0.43	10.31	41,54,69,77	20
5	BOG	A	608	20/20	0.40	8.49	48,65,100,104	20
5	BOG	B	608	20/20	0.32	7.88	28,61,78,81	20
3	NA	A	602	1/1	0.43	7.72	40,40,40,40	0
3	NA	B	604	1/1	0.36	4.53	49,49,49,49	0
5	BOG	A	606	20/20	0.34	3.96	47,79,95,100	20
5	BOG	A	609	20/20	0.71	3.43	49,69,85,88	20
5	BOG	A	610	20/20	0.48	3.29	39,84,97,97	20
2	LEU	A	601	9/9	0.24	1.38	33,42,44,45	0
3	NA	A	603	1/1	0.24	1.18	46,46,46,46	0
2	LEU	B	603	9/9	0.26	0.81	38,45,49,54	0
3	NA	B	605	1/1	0.23	0.43	55,55,55,55	0
5	BOG	B	601	20/20	0.27	0.27	44,74,92,94	20
5	BOG	A	607	20/20	0.19	0.15	50,68,80,84	0
4	BR	B	606	1/1	0.17	-0.59	77,77,77,77	1
4	BR	B	610	1/1	0.18	-1.20	88,88,88,88	1
4	BR	A	611	1/1	0.19	-1.39	76,76,76,76	1
4	BR	B	612	1/1	0.10	-1.62	90,90,90,90	1
4	BR	A	604	1/1	0.10	-1.77	73,73,73,73	1
4	BR	A	612	1/1	0.10	-2.04	96,96,96,96	1
4	BR	B	607	1/1	0.06	-2.65	75,75,75,75	1
4	BR	B	611	1/1	0.07	-2.71	84,84,84,84	1
4	BR	A	614	1/1	0.08	-2.73	63,63,63,63	1
4	BR	A	605	1/1	0.06	-3.49	59,59,59,59	1

6.5 Other polymers ⓘ

There are no such residues in this entry.