



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:17 AM UTC

PDB ID : 3JY6 / pdb_00003jy6
Title : Crystal structure of LacI Transcriptional regulator from *Lactobacillus brevis*
Authors : Syed Ibrahim, B.; Kumaran, D.; Burley, S.K.; Swaminathan, S.; New York
SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2009-09-21
Resolution : 1.97 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

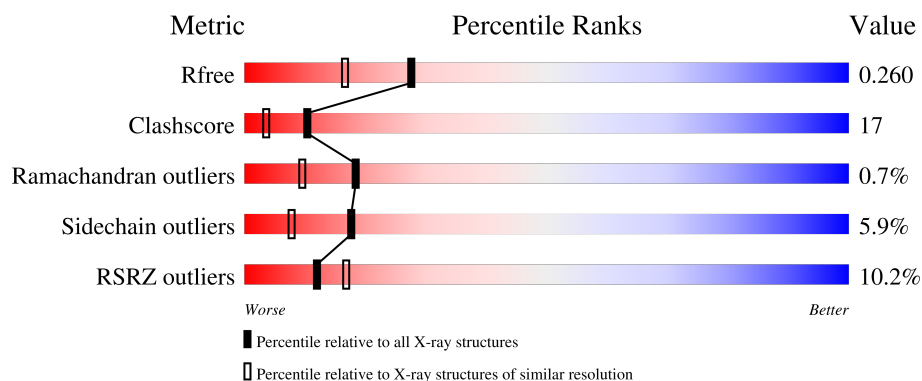
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.97 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1506 (1.98-1.98)
Clashscore	190562	1534 (1.98-1.98)
Ramachandran outliers	187476	1518 (1.98-1.98)
Sidechain outliers	187428	1518 (1.98-1.98)
RSRZ outliers	180081	1506 (1.98-1.98)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	276	3% 72% 20% • 5%
1	B	276	11% 64% 25% • • 6%
1	C	276	10% 61% 28% • • 8%
1	D	276	14% 59% 25% 9% • 5%

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	EDO	D	3318	-	-	X	-
3	EDO	D	3319	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 8291 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 Transcriptional regulator, LacI family.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2065	1307	353	398	7			
1	B	259	Total	C	N	O	S	0	0	0
			2027	1285	348	387	7			
1	C	254	Total	C	N	O	S	0	0	0
			1976	1256	337	376	7			
1	D	261	Total	C	N	O	S	0	0	0
			2038	1291	347	393	7			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	61	MET	-	expression tag	UNP Q03T70
A	62	SER	-	expression tag	UNP Q03T70
A	63	LEU	-	expression tag	UNP Q03T70
A	329	GLU	-	expression tag	UNP Q03T70
A	330	GLY	-	expression tag	UNP Q03T70
A	331	HIS	-	expression tag	UNP Q03T70
A	332	HIS	-	expression tag	UNP Q03T70
A	333	HIS	-	expression tag	UNP Q03T70
A	334	HIS	-	expression tag	UNP Q03T70
A	335	HIS	-	expression tag	UNP Q03T70
A	336	HIS	-	expression tag	UNP Q03T70
B	61	MET	-	expression tag	UNP Q03T70
B	62	SER	-	expression tag	UNP Q03T70
B	63	LEU	-	expression tag	UNP Q03T70
B	329	GLU	-	expression tag	UNP Q03T70
B	330	GLY	-	expression tag	UNP Q03T70
B	331	HIS	-	expression tag	UNP Q03T70
B	332	HIS	-	expression tag	UNP Q03T70
B	333	HIS	-	expression tag	UNP Q03T70
B	334	HIS	-	expression tag	UNP Q03T70
B	335	HIS	-	expression tag	UNP Q03T70

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Chain	Residue	Modelled	Actual	Comment	Reference
B	336	HIS	-	expression tag	UNP Q03T70
C	61	MET	-	expression tag	UNP Q03T70
C	62	SER	-	expression tag	UNP Q03T70
C	63	LEU	-	expression tag	UNP Q03T70
C	329	GLU	-	expression tag	UNP Q03T70
C	330	GLY	-	expression tag	UNP Q03T70
C	331	HIS	-	expression tag	UNP Q03T70
C	332	HIS	-	expression tag	UNP Q03T70
C	333	HIS	-	expression tag	UNP Q03T70
C	334	HIS	-	expression tag	UNP Q03T70
C	335	HIS	-	expression tag	UNP Q03T70
C	336	HIS	-	expression tag	UNP Q03T70
D	61	MET	-	expression tag	UNP Q03T70
D	62	SER	-	expression tag	UNP Q03T70
D	63	LEU	-	expression tag	UNP Q03T70
D	329	GLU	-	expression tag	UNP Q03T70
D	330	GLY	-	expression tag	UNP Q03T70
D	331	HIS	-	expression tag	UNP Q03T70
D	332	HIS	-	expression tag	UNP Q03T70
D	333	HIS	-	expression tag	UNP Q03T70
D	334	HIS	-	expression tag	UNP Q03T70
D	335	HIS	-	expression tag	UNP Q03T70
D	336	HIS	-	expression tag	UNP Q03T70

- Molecule 2 is CHLORIDE ION (CCD ID: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total Cl 1 1	0	0

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	D	1	Total	C	O	0	0
			4	2	2		
3	D	1	Total	C	O	0	0
			4	2	2		

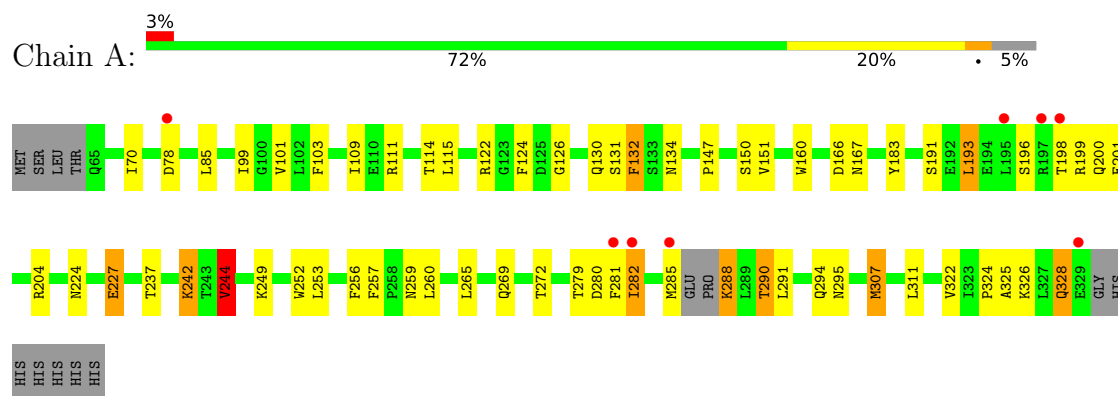
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	50	Total	O	0	0
			50	50		
4	B	34	Total	O	0	0
			34	34		
4	C	50	Total	O	0	0
			50	50		
4	D	42	Total	O	0	0
			42	42		

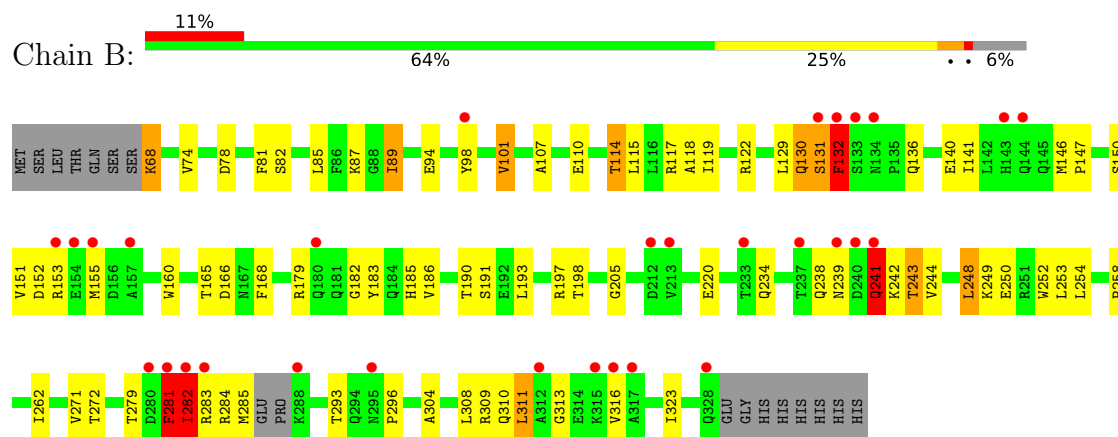
3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

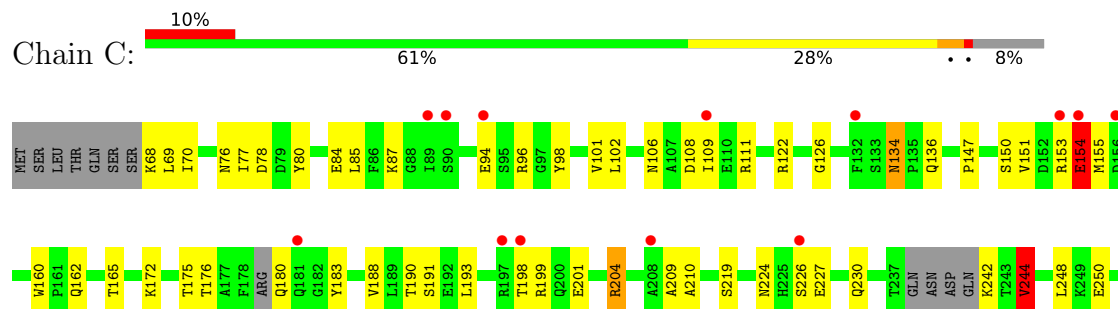
- Molecule 1: Transcriptional regulator, LacI family

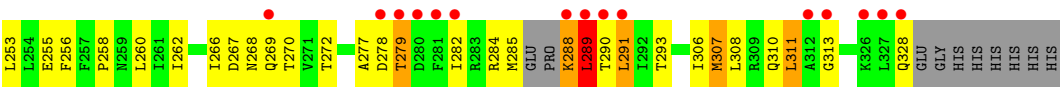


- Molecule 1: Transcriptional regulator, LacI family

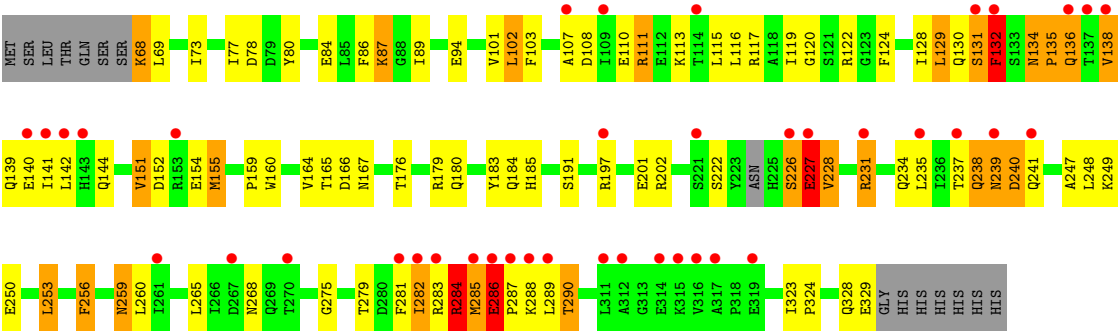


- Molecule 1: Transcriptional regulator, LacI family





● Molecule 1: Transcriptional regulator, LacI family



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	40.57Å 109.33Å 113.34Å 90.00° 95.17° 90.00°	Depositor
Resolution (Å)	40.41 – 1.97 40.41 – 1.97	Depositor EDS
% Data completeness (in resolution range)	88.3 (40.41-1.97) 88.4 (40.41-1.97)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.47 (at 1.97Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.247 , 0.261 0.247 , 0.260	Depositor DCC
R_{free} test set	3111 reflections (4.72%)	wwPDB-VP
Wilson B-factor (Å ²)	23.1	Xtriage
Anisotropy	0.221	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 36.3	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	8291	wwPDB-VP
Average B, all atoms (Å ²)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.86% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.59	1/2097 (0.0%)	1.06	14/2840 (0.5%)
1	B	0.51	0/2058	1.04	12/2788 (0.4%)
1	C	0.51	1/2005 (0.0%)	1.05	10/2714 (0.4%)
1	D	0.66	2/2070 (0.1%)	1.27	32/2807 (1.1%)
All	All	0.57	4/8230 (0.0%)	1.11	68/11149 (0.6%)

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	D	0	1

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	307	MET	SD-CE	-10.73	1.52	1.79
1	C	307	MET	SD-CE	-8.94	1.57	1.79
1	D	286	GLU	CA-CB	-6.64	1.46	1.53
1	D	129	LEU	CA-C	-5.40	1.45	1.52

All (68) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	289	LEU	N-CA-C	16.83	131.93	111.40
1	B	282	ILE	CB-CA-C	-11.18	97.27	112.24
1	D	286	GLU	N-CA-C	-10.73	89.35	108.55
1	D	286	GLU	O-C-N	10.72	128.73	121.23
1	D	132	PHE	N-CA-C	9.77	131.62	110.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	139	GLN	N-CA-C	9.54	124.02	112.38
1	D	238	GLN	N-CA-C	-9.11	101.33	111.07
1	D	226	SER	N-CA-C	-8.06	101.58	111.33
1	A	282	ILE	CB-CA-C	-8.00	99.12	112.16
1	A	191	SER	N-CA-C	-7.73	100.11	110.55
1	B	241	GLN	N-CA-C	-7.67	99.91	110.35
1	D	284	ARG	N-CA-C	-7.58	94.65	110.80
1	B	131	SER	N-CA-C	7.44	117.82	108.45
1	A	288	LYS	CA-C-O	7.40	133.38	120.80
1	D	239	ASN	N-CA-C	-7.36	99.36	109.95
1	D	286	GLU	N-CA-CB	7.15	121.99	111.77
1	B	129	LEU	N-CA-C	7.13	120.03	108.41
1	C	191	SER	N-CA-C	-7.12	101.34	110.53
1	C	282	ILE	N-CA-C	-7.11	104.04	111.58
1	D	134	ASN	N-CA-C	-6.99	99.40	110.10
1	D	289	LEU	N-CA-C	-6.77	98.22	109.46
1	B	193	LEU	N-CA-C	6.62	118.15	111.07
1	D	286	GLU	CB-CA-C	-6.54	102.50	110.81
1	A	132	PHE	N-CA-C	6.39	121.27	112.90
1	D	138	VAL	N-CA-C	-6.35	104.32	110.42
1	D	286	GLU	CA-C-N	6.34	127.77	119.84
1	D	286	GLU	C-N-CA	6.34	127.77	119.84
1	D	129	LEU	N-CA-C	6.26	119.70	109.24
1	D	191	SER	N-CA-C	-6.26	102.46	110.53
1	C	154	GLU	N-CA-C	-6.21	96.19	107.75
1	D	167	ASN	N-CA-C	6.20	118.58	111.02
1	D	124	PHE	N-CA-C	-6.04	102.39	110.55
1	D	323	ILE	N-CA-C	-5.98	101.20	107.60
1	C	269	GLN	N-CA-C	-5.96	105.00	112.93
1	B	191	SER	N-CA-C	-5.94	102.87	110.53
1	A	291	LEU	N-CA-C	5.93	119.22	109.85
1	A	124	PHE	N-CA-C	-5.91	100.65	110.17
1	B	132	PHE	N-CA-C	5.90	120.60	113.17
1	B	250	GLU	N-CA-C	5.87	118.43	111.33
1	D	134	ASN	CA-C-N	5.87	127.18	119.84
1	D	134	ASN	C-N-CA	5.87	127.18	119.84
1	D	130	GLN	N-CA-C	-5.86	104.27	112.12
1	D	129	LEU	CA-CB-CG	5.82	136.68	116.30
1	C	244	VAL	CB-CA-C	-5.82	102.51	110.77
1	A	244	VAL	CB-CA-C	-5.69	102.69	110.77
1	D	131	SER	N-CA-C	5.55	122.63	110.80
1	A	288	LYS	N-CA-CB	5.55	119.93	110.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	328	GLN	N-CA-C	5.55	117.45	108.41
1	D	290	THR	N-CA-C	-5.55	100.67	109.76
1	A	256	PHE	N-CA-C	5.52	121.04	113.97
1	A	242	LYS	N-CA-C	-5.41	102.38	110.23
1	D	323	ILE	CB-CA-C	-5.41	104.95	111.18
1	C	193	LEU	N-CA-C	5.39	116.84	111.07
1	B	281	PHE	N-CA-C	5.38	119.83	113.16
1	A	193	LEU	N-CA-C	5.36	116.80	111.07
1	D	102	LEU	N-CA-C	5.33	117.93	109.24
1	D	286	GLU	CB-CG-CD	-5.27	103.64	112.60
1	D	222	SER	N-CA-C	5.26	116.77	107.61
1	A	167	ASN	N-CA-C	5.26	117.42	111.11
1	D	256	PHE	N-CA-C	5.23	120.66	113.97
1	A	237	THR	N-CA-C	5.20	119.60	113.16
1	C	313	GLY	N-CA-C	5.18	122.43	115.59
1	B	146	MET	N-CA-C	5.11	114.17	108.25
1	B	243	THR	N-CA-C	5.11	118.07	109.95
1	C	311	LEU	N-CA-C	-5.09	106.58	112.89
1	B	323	ILE	N-CA-C	-5.08	102.94	107.56
1	C	291	LEU	N-CA-C	5.05	117.54	109.81
1	D	227	GLU	N-CA-C	-5.04	101.09	109.46

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	D	286	GLU	Mainchain

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	2065	0	2066	51	0
1	B	2027	0	2028	73	0
1	C	1976	0	1978	79	0
1	D	2038	0	2015	98	0
2	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	8	0	12	16	0
4	A	50	0	0	5	0
4	B	34	0	0	5	0
4	C	50	0	0	9	0
4	D	42	0	0	6	0
All	All	8291	0	8099	275	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 17.

All (275) 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:154:GLU:OE1	1:C:201:GLU:HG3	1.51	1.09
1:A:134:ASN:HB2	4:A:355:HOH:O	1.64	0.97
1:C:68:LYS:HD3	4:C:379:HOH:O	1.65	0.97
1:D:107:ALA:HA	1:D:132:PHE:HZ	1.33	0.92
1:C:134:ASN:HD21	1:C:136:GLN:HB2	1.32	0.91
1:D:107:ALA:HA	1:D:132:PHE:CZ	2.05	0.90
1:A:70:ILE:HD12	1:A:307:MET:HE2	1.59	0.85
1:B:283:ARG:HH22	1:B:293:THR:CG2	1.93	0.81
1:B:185:HIS:HB2	1:B:241:GLN:HG3	1.62	0.81
3:D:3318:EDO:H12	4:D:4:HOH:O	1.80	0.80
1:C:94:GLU:OE2	1:D:111:ARG:HD2	1.83	0.79
1:B:283:ARG:NH2	1:B:293:THR:CG2	2.47	0.78
1:D:134:ASN:HD21	1:D:136:GLN:HG3	1.48	0.78
1:C:224:ASN:HD22	1:C:227:GLU:H	1.31	0.77
1:C:147:PRO:HG2	1:C:307:MET:HE1	1.66	0.77
1:A:122:ARG:HH21	1:B:68:LYS:N	1.84	0.76
1:D:154:GLU:HG3	1:D:164:VAL:HG11	1.68	0.76
1:D:279:THR:HB	1:D:282:ILE:CB	2.15	0.75
1:A:224:ASN:HB3	1:A:227:GLU:HB3	1.70	0.74
1:C:109:ILE:HD11	4:C:346:HOH:O	1.87	0.74
1:D:281:PHE:CD1	1:D:281:PHE:C	2.65	0.73
1:A:285:MET:HE3	1:B:258:PRO:HB3	1.70	0.73
1:C:255:GLU:HG3	1:D:284:ARG:HH12	1.54	0.73
3:D:3318:EDO:H21	3:D:3319:EDO:H22	1.71	0.72
1:C:126:GLY:HA3	1:C:307:MET:HE2	1.71	0.72
3:D:3318:EDO:H21	3:D:3319:EDO:C2	2.19	0.72
1:D:239:ASN:ND2	1:D:240:ASP:H	1.88	0.72
1:B:74:VAL:HA	1:B:130:GLN:HG3	1.70	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:328:GLN:HB3	4:C:380:HOH:O	1.90	0.71
1:C:183:TYR:HB3	1:C:244:VAL:HG22	1.72	0.71
1:A:147:PRO:HG2	1:A:307:MET:HE1	1.73	0.71
1:D:283:ARG:O	1:D:284:ARG:HG3	1.90	0.71
1:D:227:GLU:HB3	1:D:231:ARG:HH12	1.55	0.71
1:A:279:THR:HG21	1:A:281:PHE:CZ	2.25	0.70
1:A:285:MET:HE2	1:B:262:ILE:HD11	1.73	0.70
1:A:78:ASP:OD2	1:B:87:LYS:HE3	1.91	0.70
1:B:85:LEU:HD11	1:B:151:VAL:HG11	1.74	0.69
1:C:76:ASN:OD1	1:C:78:ASP:HB2	1.92	0.69
1:C:94:GLU:HG3	1:D:111:ARG:NH1	2.06	0.69
1:B:279:THR:HG21	1:B:281:PHE:CE2	2.29	0.68
1:A:279:THR:HG21	1:A:281:PHE:CE2	2.29	0.68
1:A:249:LYS:HG2	4:A:24:HOH:O	1.92	0.68
1:C:85:LEU:HD11	1:C:151:VAL:HG11	1.76	0.67
1:D:135:PRO:CD	1:D:155:MET:HE3	2.24	0.67
1:A:126:GLY:HA3	1:A:307:MET:CE	2.25	0.67
1:B:185:HIS:CB	1:B:241:GLN:HG3	2.24	0.66
1:A:126:GLY:HA3	1:A:307:MET:HE2	1.77	0.66
1:C:199:ARG:HG2	4:C:347:HOH:O	1.96	0.66
1:D:134:ASN:HD21	1:D:136:GLN:CG	2.08	0.66
1:C:80:TYR:CZ	1:C:84:GLU:HG3	2.31	0.65
1:C:102:LEU:HD23	1:D:102:LEU:HD23	1.79	0.65
1:C:183:TYR:HB3	1:C:244:VAL:CG2	2.27	0.65
1:A:183:TYR:HB3	1:A:244:VAL:HG22	1.78	0.65
1:C:262:ILE:HD12	1:D:285:MET:HG2	1.78	0.65
1:D:250:GLU:H	3:D:3318:EDO:C2	2.09	0.65
1:B:253:LEU:CD2	1:B:282:ILE:HG21	2.27	0.64
1:C:126:GLY:HA3	1:C:307:MET:CE	2.27	0.64
1:C:224:ASN:ND2	1:C:227:GLU:H	1.96	0.64
1:D:249:LYS:HG2	3:D:3318:EDO:H22	1.80	0.64
1:A:259:ASN:HD21	1:B:285:MET:HA	1.63	0.64
3:D:3318:EDO:H21	3:D:3319:EDO:O2	1.97	0.64
1:C:224:ASN:ND2	1:C:227:GLU:HG3	2.13	0.63
1:D:120:GLY:HA3	1:D:142:LEU:O	1.98	0.63
1:B:279:THR:HG22	1:B:281:PHE:H	1.63	0.63
1:A:259:ASN:ND2	1:B:285:MET:HA	2.14	0.62
1:A:198:THR:HG23	4:A:343:HOH:O	2.00	0.62
1:D:197:ARG:O	1:D:201:GLU:HG3	1.99	0.62
1:D:283:ARG:C	1:D:284:ARG:HG3	2.25	0.62
1:A:307:MET:HE3	1:A:311:LEU:HG	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:76:ASN:HD21	1:D:87:LYS:NZ	1.97	0.61
1:D:68:LYS:NZ	1:D:68:LYS:HB3	2.14	0.61
1:D:249:LYS:HA	3:D:3318:EDO:H22	1.83	0.61
1:A:99:ILE:HD12	1:B:122:ARG:HH12	1.66	0.61
1:D:281:PHE:O	1:D:283:ARG:N	2.34	0.60
1:C:165:THR:HG23	4:C:342:HOH:O	2.01	0.60
1:C:262:ILE:CD1	1:D:285:MET:HG2	2.31	0.60
1:D:176:THR:O	1:D:180:GLN:HG3	2.02	0.60
1:A:307:MET:HE3	1:A:311:LEU:CG	2.32	0.60
1:B:253:LEU:HD21	1:B:282:ILE:HG21	1.84	0.59
1:D:281:PHE:C	1:D:283:ARG:H	2.11	0.59
1:B:283:ARG:NH2	1:B:293:THR:HG21	2.18	0.59
1:A:279:THR:CG2	1:A:281:PHE:CE2	2.85	0.59
1:D:227:GLU:CD	1:D:227:GLU:H	2.11	0.58
1:A:242:LYS:HE3	1:A:269:GLN:O	2.04	0.58
1:C:96:ARG:HG3	1:C:308:LEU:HD13	1.85	0.58
1:C:190:THR:HB	1:C:248:LEU:HD22	1.85	0.58
1:D:249:LYS:HG2	3:D:3318:EDO:C1	2.33	0.58
1:A:101:VAL:HG13	1:B:101:VAL:HG13	1.85	0.58
1:C:154:GLU:OE1	1:C:201:GLU:CG	2.39	0.58
1:D:253:LEU:HD11	1:D:275:GLY:HA3	1.84	0.58
1:B:258:PRO:HG3	1:B:285:MET:HE1	1.85	0.57
1:D:288:LYS:HA	4:D:355:HOH:O	2.03	0.57
1:D:281:PHE:C	1:D:283:ARG:N	2.62	0.57
1:C:250:GLU:OE2	1:C:277:ALA:HB1	2.04	0.56
1:A:244:VAL:HG13	1:A:272:THR:HG23	1.86	0.56
1:B:110:GLU:O	1:B:114:THR:HG22	2.04	0.56
1:B:281:PHE:CD1	1:B:281:PHE:C	2.84	0.56
1:D:202:ARG:HH12	3:D:3319:EDO:H21	1.69	0.56
1:D:249:LYS:HG2	3:D:3318:EDO:H12	1.87	0.56
1:C:307:MET:HE3	1:C:311:LEU:CG	2.35	0.56
1:D:286:GLU:C	1:D:288:LYS:H	2.13	0.56
1:B:190:THR:HB	1:B:248:LEU:HD22	1.88	0.56
1:C:70:ILE:HD12	1:C:307:MET:HE2	1.87	0.55
1:B:136:GLN:O	1:B:140:GLU:HG3	2.06	0.55
1:D:135:PRO:HD3	1:D:155:MET:HE3	1.86	0.55
1:B:197:ARG:CB	4:B:351:HOH:O	2.54	0.55
1:C:153:ARG:HB2	4:C:343:HOH:O	2.07	0.55
1:D:290:THR:CG2	1:D:329:GLU:HG3	2.37	0.55
1:B:283:ARG:NH2	1:B:293:THR:HG22	2.21	0.55
1:B:131:SER:O	1:B:153:ARG:HG2	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:198:THR:HG23	4:C:353:HOH:O	2.05	0.54
1:C:328:GLN:HG2	4:C:380:HOH:O	2.07	0.54
1:D:113:LYS:NZ	1:D:141:ILE:HD11	2.21	0.54
1:B:283:ARG:NH1	4:B:352:HOH:O	2.39	0.54
1:D:231:ARG:HD2	4:D:360:HOH:O	2.07	0.54
1:A:253:LEU:HG	1:A:257:PHE:CD2	2.43	0.54
1:D:239:ASN:OD1	1:D:241:GLN:HB2	2.08	0.54
1:C:307:MET:HE3	1:C:311:LEU:HD11	1.89	0.54
1:A:166:ASP:OD1	1:A:324:PRO:HA	2.08	0.53
1:C:188:VAL:HG13	1:C:248:LEU:HD13	1.89	0.53
1:C:328:GLN:HG3	1:C:328:GLN:O	2.09	0.52
1:C:154:GLU:HG3	1:C:198:THR:HG22	1.90	0.52
1:C:284:ARG:HB3	1:C:284:ARG:NH1	2.25	0.52
1:B:239:ASN:CB	1:B:241:GLN:CD	2.83	0.52
1:B:81:PHE:CD1	1:B:296:PRO:HB3	2.44	0.52
1:B:115:LEU:O	1:B:119:ILE:HD13	2.08	0.52
1:C:154:GLU:CG	1:C:198:THR:HG22	2.40	0.52
1:C:101:VAL:CG1	1:D:101:VAL:HG13	2.40	0.52
1:C:108:ASP:HB3	1:C:111:ARG:HB3	1.91	0.51
1:D:279:THR:HG22	1:D:282:ILE:H	1.76	0.51
1:C:253:LEU:O	1:C:258:PRO:HD3	2.10	0.51
1:C:68:LYS:HB3	1:C:98:TYR:CD1	2.46	0.51
1:D:111:ARG:O	1:D:115:LEU:HG	2.11	0.51
1:C:85:LEU:CD1	1:C:151:VAL:HG11	2.41	0.51
1:B:279:THR:HG21	1:B:281:PHE:CD2	2.46	0.51
1:B:85:LEU:O	1:B:89:ILE:HD12	2.11	0.51
1:A:109:ILE:HD13	1:A:132:PHE:HB2	1.94	0.50
1:C:256:PHE:O	1:C:260:LEU:HD13	2.10	0.50
1:B:238:GLN:HG3	1:B:239:ASN:N	2.26	0.50
1:B:150:SER:HB3	1:B:160:TRP:HE1	1.76	0.50
1:B:118:ALA:HB1	1:B:122:ARG:HH11	1.76	0.50
1:D:249:LYS:HG2	3:D:3318:EDO:C2	2.41	0.50
1:D:108:ASP:OD1	1:D:110:GLU:HB2	2.11	0.50
1:D:290:THR:HG21	1:D:329:GLU:HG3	1.94	0.50
1:B:183:TYR:CD1	1:B:242:LYS:HG2	2.47	0.49
1:B:239:ASN:CB	1:B:241:GLN:NE2	2.75	0.49
1:D:103:PHE:HB3	1:D:115:LEU:HD13	1.94	0.49
1:D:151:VAL:HG12	1:D:152:ASP:N	2.27	0.49
1:C:307:MET:HE3	1:C:311:LEU:HG	1.94	0.49
1:C:176:THR:O	1:C:180:GLN:HG3	2.12	0.49
1:C:224:ASN:HD22	1:C:227:GLU:N	2.04	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:185:HIS:HE1	1:D:235:LEU:O	1.96	0.49
1:B:165:THR:HG22	1:B:166:ASP:N	2.27	0.49
1:B:197:ARG:HB2	4:B:351:HOH:O	2.12	0.49
1:C:155:MET:CB	1:C:162:GLN:HE22	2.26	0.48
1:A:244:VAL:HG13	1:A:272:THR:CG2	2.44	0.48
1:C:277:ALA:HB3	1:C:291:LEU:HD22	1.95	0.48
1:A:259:ASN:ND2	1:B:284:ARG:O	2.44	0.48
1:C:76:ASN:HD21	1:D:87:LYS:HZ2	1.62	0.48
1:D:260:LEU:HD22	1:D:265:LEU:HD23	1.96	0.48
1:C:267:ASP:O	1:C:268:ASN:HB2	2.14	0.48
1:B:117:ARG:HG2	1:B:141:ILE:HG23	1.95	0.47
1:A:85:LEU:HD22	1:A:130:GLN:HG3	1.96	0.47
1:B:249:LYS:HG2	4:B:337:HOH:O	2.13	0.47
1:B:279:THR:CG2	1:B:281:PHE:CD2	2.98	0.47
1:D:113:LYS:O	1:D:117:ARG:HG3	2.14	0.47
1:A:294:GLN:O	1:A:295:ASN:HB2	2.15	0.47
1:C:87:LYS:HG3	1:D:77:ILE:HD11	1.95	0.47
1:A:285:MET:HE3	1:B:258:PRO:CB	2.40	0.47
1:B:220:GLU:HG2	1:B:252:TRP:HZ2	1.79	0.47
1:D:239:ASN:ND2	1:D:240:ASP:N	2.60	0.47
1:D:290:THR:HB	1:D:329:GLU:HG3	1.97	0.47
1:C:77:ILE:HD13	1:D:86:PHE:HD2	1.81	0.46
1:D:80:TYR:CZ	1:D:84:GLU:HG3	2.51	0.46
1:D:250:GLU:H	3:D:3318:EDO:H22	1.78	0.46
1:A:260:LEU:HG	1:A:265:LEU:HD23	1.97	0.46
1:D:247:ALA:HB3	1:D:253:LEU:CD1	2.46	0.46
1:D:249:LYS:CG	3:D:3318:EDO:H22	2.45	0.46
1:B:152:ASP:O	1:B:198:THR:HG21	2.16	0.46
1:C:307:MET:HE3	1:C:311:LEU:CD1	2.46	0.46
1:C:201:GLU:OE1	1:C:204:ARG:NH1	2.49	0.46
1:C:328:GLN:CG	4:C:380:HOH:O	2.62	0.45
1:D:68:LYS:HB3	1:D:68:LYS:HZ3	1.80	0.45
1:A:111:ARG:NH1	1:B:94:GLU:HG3	2.31	0.45
1:B:238:GLN:HG3	1:B:239:ASN:H	1.81	0.45
1:C:111:ARG:NH1	1:D:94:GLU:HG3	2.32	0.45
1:C:172:LYS:HG3	1:C:209:ALA:HB2	1.98	0.45
1:B:82:SER:OG	1:B:130:GLN:NE2	2.49	0.45
1:B:183:TYR:N	1:B:242:LYS:HZ2	2.14	0.45
3:D:3318:EDO:C2	3:D:3319:EDO:H22	2.43	0.45
1:D:165:THR:HG22	1:D:166:ASP:N	2.32	0.45
1:D:227:GLU:O	1:D:228:VAL:CB	2.64	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:168:PHE:CE1	1:B:205:GLY:HA2	2.52	0.45
1:A:322:VAL:HG12	4:A:10:HOH:O	2.16	0.45
1:C:226:SER:O	1:C:230:GLN:HG3	2.17	0.45
1:C:69:LEU:HD11	1:D:122:ARG:HD3	1.98	0.45
1:D:166:ASP:OD1	1:D:324:PRO:HA	2.16	0.45
1:D:159:PRO:HG2	1:D:160:TRP:CE3	2.52	0.44
1:D:227:GLU:O	1:D:228:VAL:HG23	2.17	0.44
1:C:101:VAL:HG13	1:D:101:VAL:HG13	1.99	0.44
1:C:106:ASN:OD1	1:D:87:LYS:NZ	2.50	0.44
1:C:242:LYS:HA	1:C:270:THR:O	2.18	0.44
1:C:278:ASP:OD1	1:C:293:THR:HB	2.17	0.44
1:D:110:GLU:OE1	1:D:113:LYS:HD2	2.18	0.44
1:B:304:ALA:O	1:B:308:LEU:HG	2.18	0.44
1:D:250:GLU:CB	3:D:3318:EDO:O2	2.65	0.44
1:C:289:LEU:O	1:C:290:THR:HB	2.16	0.44
1:D:135:PRO:CG	1:D:155:MET:HE3	2.48	0.44
1:C:150:SER:HB2	1:C:160:TRP:HE1	1.83	0.44
1:D:281:PHE:CD1	1:D:282:ILE:N	2.86	0.44
1:D:286:GLU:C	1:D:288:LYS:N	2.76	0.44
1:A:103:PHE:HB3	1:A:115:LEU:HD13	2.00	0.43
1:A:259:ASN:HD21	1:B:284:ARG:C	2.25	0.43
1:B:107:ALA:HA	1:B:132:PHE:CZ	2.53	0.43
1:C:134:ASN:ND2	1:C:136:GLN:HB2	2.15	0.43
1:C:224:ASN:HD22	1:C:227:GLU:HG3	1.84	0.43
1:B:147:PRO:HB3	1:B:310:GLN:HB3	2.00	0.43
1:A:183:TYR:CB	1:A:244:VAL:HG22	2.48	0.43
1:D:328:GLN:NE2	4:D:354:HOH:O	2.44	0.43
1:D:135:PRO:HG3	1:D:155:MET:HE3	2.01	0.43
1:B:185:HIS:CG	1:B:241:GLN:HG3	2.54	0.42
1:A:78:ASP:CG	1:B:87:LYS:HE3	2.44	0.42
1:A:328:GLN:O	1:A:328:GLN:HG3	2.19	0.42
1:B:130:GLN:HG3	1:B:130:GLN:O	2.19	0.42
1:B:182:GLY:C	1:B:242:LYS:HZ2	2.27	0.42
1:D:281:PHE:C	1:D:281:PHE:HD1	2.22	0.42
1:D:227:GLU:O	1:D:228:VAL:HB	2.18	0.42
1:D:259:ASN:HD22	1:D:259:ASN:HA	1.64	0.42
1:B:220:GLU:HG2	1:B:252:TRP:CZ2	2.54	0.42
1:C:175:THR:CG2	1:C:210:ALA:HB2	2.49	0.42
1:A:193:LEU:O	1:A:200:GLN:HG3	2.20	0.42
1:A:325:ALA:O	1:A:326:LYS:HD3	2.20	0.42
1:B:186:VAL:HG22	1:B:244:VAL:CG1	2.50	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:266:ILE:HG23	1:C:272:THR:HA	2.01	0.42
1:D:285:MET:HE3	1:D:285:MET:HB2	1.63	0.42
1:D:116:LEU:HD22	1:D:142:LEU:HD13	2.01	0.42
1:B:241:GLN:HE21	1:B:241:GLN:HB3	1.51	0.42
1:A:201:GLU:OE2	1:A:204:ARG:NH1	2.53	0.42
1:D:179:ARG:HA	1:D:183:TYR:O	2.20	0.42
1:A:196:SER:HB3	1:A:199:ARG:CD	2.50	0.42
1:A:285:MET:HE2	1:B:262:ILE:CD1	2.48	0.41
1:B:244:VAL:HG23	1:B:272:THR:HG23	2.01	0.41
1:A:85:LEU:HD11	1:A:151:VAL:HG11	2.03	0.41
1:A:126:GLY:HA3	1:A:307:MET:HE1	1.99	0.41
1:A:150:SER:HB2	1:A:160:TRP:HE1	1.85	0.41
1:C:250:GLU:OE2	1:C:279:THR:HB	2.20	0.41
1:D:250:GLU:HB2	3:D:3318:EDO:O2	2.19	0.41
1:A:280:ASP:HB2	4:A:338:HOH:O	2.20	0.41
1:B:254:LEU:HD21	1:B:282:ILE:HD13	2.01	0.41
1:C:87:LYS:HE3	1:D:78:ASP:OD2	2.19	0.41
1:C:122:ARG:HE	1:D:69:LEU:HD12	1.85	0.41
1:B:68:LYS:HB2	1:B:98:TYR:CD1	2.54	0.41
1:C:306:ILE:O	1:C:310:GLN:HG3	2.21	0.41
1:B:197:ARG:HB3	4:B:351:HOH:O	2.17	0.41
1:B:243:THR:HG22	1:B:244:VAL:N	2.36	0.41
1:D:111:ARG:NH2	4:D:345:HOH:O	2.54	0.41
1:D:113:LYS:HZ2	1:D:141:ILE:HD11	1.85	0.41
1:A:290:THR:O	1:A:290:THR:HG22	2.21	0.41
1:B:311:LEU:C	1:B:313:GLY:H	2.28	0.41
1:C:77:ILE:HG22	1:C:102:LEU:HD21	2.03	0.41
1:D:73:ILE:HD11	1:D:119:ILE:HD12	2.03	0.41
1:D:89:ILE:CD1	1:D:128:ILE:HG21	2.51	0.41
1:D:231:ARG:CD	4:D:360:HOH:O	2.68	0.41
1:C:284:ARG:HG3	1:C:284:ARG:O	2.21	0.41
1:D:234:GLN:O	1:D:238:GLN:N	2.53	0.41
1:D:290:THR:HG21	1:D:329:GLU:CG	2.50	0.41
1:D:290:THR:CB	1:D:329:GLU:HA	2.51	0.41
1:B:183:TYR:CG	1:B:242:LYS:NZ	2.89	0.40
1:C:288:LYS:O	1:C:288:LYS:HG2	2.20	0.40
1:D:290:THR:HB	1:D:329:GLU:HA	2.02	0.40
1:A:252:TRP:CE3	1:A:252:TRP:HA	2.57	0.40
1:B:309:ARG:NH1	1:B:316:VAL:HG13	2.37	0.40
1:D:256:PHE:O	1:D:260:LEU:HG	2.21	0.40
1:B:243:THR:O	1:B:271:VAL:HA	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:138:VAL:O	1:D:142:LEU:HB2	2.22	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	259/276 (94%)	252 (97%)	7 (3%)	0	100	100
1	B	255/276 (92%)	247 (97%)	8 (3%)	0	100	100
1	C	246/276 (89%)	233 (95%)	13 (5%)	0	100	100
1	D	257/276 (93%)	242 (94%)	8 (3%)	7 (3%)	4	0
All	All	1017/1104 (92%)	974 (96%)	36 (4%)	7 (1%)	18	9

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	131	SER
1	D	228	VAL
1	D	282	ILE
1	D	287	PRO
1	D	132	PHE
1	D	284	ARG
1	D	285	MET

5.3.2 Protein sidechains [i](#)

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	226/240 (94%)	219 (97%)	7 (3%)	35	26
1	B	219/240 (91%)	204 (93%)	15 (7%)	14	5
1	C	213/240 (89%)	204 (96%)	9 (4%)	26	15
1	D	219/240 (91%)	198 (90%)	21 (10%)	8	2
All	All	877/960 (91%)	825 (94%)	52 (6%)	18	8

All (52) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	114	THR
1	A	131	SER
1	A	227	GLU
1	A	244	VAL
1	A	282	ILE
1	A	288	LYS
1	A	290	THR
1	B	68	LYS
1	B	78	ASP
1	B	89	ILE
1	B	101	VAL
1	B	114	THR
1	B	130	GLN
1	B	132	PHE
1	B	155	MET
1	B	179	ARG
1	B	234	GLN
1	B	241	GLN
1	B	248	LEU
1	B	281	PHE
1	B	282	ILE
1	B	311	LEU
1	C	134	ASN
1	C	154	GLU
1	C	204	ARG
1	C	219	SER
1	C	244	VAL
1	C	279	THR
1	C	285	MET
1	C	288	LYS

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Mol	Chain	Res	Type
1	C	289	LEU
1	D	68	LYS
1	D	87	LYS
1	D	111	ARG
1	D	129	LEU
1	D	132	PHE
1	D	135	PRO
1	D	136	GLN
1	D	140	GLU
1	D	144	GLN
1	D	151	VAL
1	D	155	MET
1	D	184	GLN
1	D	226	SER
1	D	227	GLU
1	D	231	ARG
1	D	237	THR
1	D	240	ASP
1	D	248	LEU
1	D	253	LEU
1	D	259	ASN
1	D	268	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	143	HIS
1	A	144	GLN
1	A	162	GLN
1	A	180	GLN
1	A	224	ASN
1	A	230	GLN
1	A	234	GLN
1	A	259	ASN
1	A	295	ASN
1	B	130	GLN
1	B	145	GLN
1	B	162	GLN
1	B	184	GLN
1	B	295	ASN
1	B	328	GLN
1	C	76	ASN

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Mol	Chain	Res	Type
1	C	134	ASN
1	C	136	GLN
1	C	143	HIS
1	C	162	GLN
1	C	224	ASN
1	C	229	HIS
1	C	230	GLN
1	C	259	ASN
1	D	134	ASN
1	D	139	GLN
1	D	184	GLN
1	D	239	ASN
1	D	241	GLN
1	D	259	ASN
1	D	295	ASN
1	D	310	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 oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 3 ligands modelled in this entry, 1 is monoatomic - leaving 2 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
3	EDO	D	3319	-	3,3,3	0.43	0	2,2,2	1.23	0
3	EDO	D	3318	-	3,3,3	0.63	0	2,2,2	0.35	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	EDO	D	3319	-	-	0/1/1/1	-
3	EDO	D	3318	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	D	3318	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 16 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	D	3319	EDO	5	0
3	D	3318	EDO	15	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	263/276 (95%)	0.48	8 (3%) 52 62	11, 24, 40, 47	1 (0%)
1	B	259/276 (93%)	0.87	30 (11%) 9 13	11, 28, 46, 51	3 (1%)
1	C	254/276 (92%)	0.83	28 (11%) 10 14	11, 29, 44, 52	1 (0%)
1	D	261/276 (94%)	0.93	40 (15%) 5 7	7, 28, 47, 58	3 (1%)
All	All	1037/1104 (93%)	0.78	106 (10%) 12 17	7, 27, 45, 58	8 (0%)

All (106) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	240	ASP	8.4
1	B	239	ASN	7.5
1	B	281	PHE	5.8
1	C	281	PHE	4.8
1	C	289	LEU	4.6
1	D	143	HIS	4.2
1	D	142	LEU	4.1
1	B	132	PHE	4.0
1	D	312	ALA	4.0
1	B	133	SER	4.0
1	C	282	ILE	3.8
1	D	315	LYS	3.8
1	D	314	GLU	3.6
1	D	282	ILE	3.6
1	A	282	ILE	3.5
1	C	327	LEU	3.4
1	D	287	PRO	3.3
1	D	221	SER	3.3
1	D	281	PHE	3.2
1	B	241	GLN	3.2
1	D	141	ILE	3.2

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Mol	Chain	Res	Type	RSRZ
1	D	197	ARG	3.2
1	D	136	GLN	3.1
1	B	316	VAL	3.1
1	D	131	SER	3.1
1	D	226	SER	3.1
1	D	114	THR	3.1
1	B	328	GLN	3.0
1	D	138	VAL	3.0
1	A	285	MET	3.0
1	B	143	HIS	3.0
1	C	312	ALA	3.0
1	D	289	LEU	3.0
1	C	288	LYS	3.0
1	B	282	ILE	3.0
1	D	237	THR	2.9
1	A	281	PHE	2.9
1	D	132	PHE	2.9
1	D	317	ALA	2.8
1	C	132	PHE	2.8
1	C	181	GLN	2.8
1	B	288	LYS	2.8
1	C	156	ASP	2.7
1	C	291	LEU	2.7
1	D	286	GLU	2.7
1	C	208	ALA	2.7
1	B	280	ASP	2.6
1	B	153	ARG	2.6
1	D	137	THR	2.6
1	B	155	MET	2.6
1	D	285	MET	2.6
1	B	154	GLU	2.6
1	D	288	LYS	2.6
1	C	153	ARG	2.5
1	B	212	ASP	2.4
1	B	295	ASN	2.4
1	C	94	GLU	2.4
1	D	267	ASP	2.4
1	C	290	THR	2.4
1	C	269	GLN	2.4
1	D	239	ASN	2.4
1	C	328	GLN	2.4
1	A	78	ASP	2.3

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Mol	Chain	Res	Type	RSRZ
1	B	283	ARG	2.3
1	D	261	ILE	2.3
1	A	195	LEU	2.3
1	D	227	GLU	2.3
1	D	241	GLN	2.3
1	B	98	TYR	2.3
1	C	197	ARG	2.3
1	C	326	LYS	2.3
1	C	313	GLY	2.2
1	D	107	ALA	2.2
1	B	131	SER	2.2
1	A	329	GLU	2.2
1	C	90	SER	2.2
1	B	157	ALA	2.2
1	B	315	LYS	2.2
1	D	235	LEU	2.2
1	C	280	ASP	2.2
1	B	144	GLN	2.2
1	A	198	THR	2.2
1	C	278	ASP	2.1
1	D	153	ARG	2.1
1	D	231	ARG	2.1
1	D	283	ARG	2.1
1	D	140	GLU	2.1
1	D	109	ILE	2.1
1	C	198	THR	2.1
1	C	154	GLU	2.1
1	D	311	LEU	2.1
1	B	180	GLN	2.1
1	C	109	ILE	2.1
1	A	197	ARG	2.1
1	C	226	SER	2.1
1	C	279	THR	2.1
1	B	213	VAL	2.1
1	B	317	ALA	2.1
1	B	134	ASN	2.0
1	B	312	ALA	2.0
1	D	319	GLU	2.0
1	C	89	ILE	2.0
1	B	233	THR	2.0
1	B	237	THR	2.0
1	D	270	THR	2.0

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Mol	Chain	Res	Type	RSRZ
1	D	316	VAL	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 oligosaccharides 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. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
3	EDO	D	3318	4/4	0.75	0.32	31,31,32,33	0
3	EDO	D	3319	4/4	0.75	0.18	38,39,40,40	0
2	CL	D	1	1/1	0.96	0.07	27,27,27,27	0

6.5 Other polymers [i](#)

There are no such residues in this entry.