



Full wwPDB X-ray Structure Validation Report ⓘ

Apr 26, 2026 – 05:47 AM EDT

PDB ID : 7OCJ / pdb_00007ocj
Title : Crystal structure of E.coli LexA in complex with nanobody NbSOS2(Nb14509)
Authors : Maso, L.; Vascon, F.; Chinellato, M.; Pardon, E.; Steyaert, J.; Angelini, A.;
Tondi, D.; Cendron, L.
Deposited on : 2021-04-27
Resolution : 2.70 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:

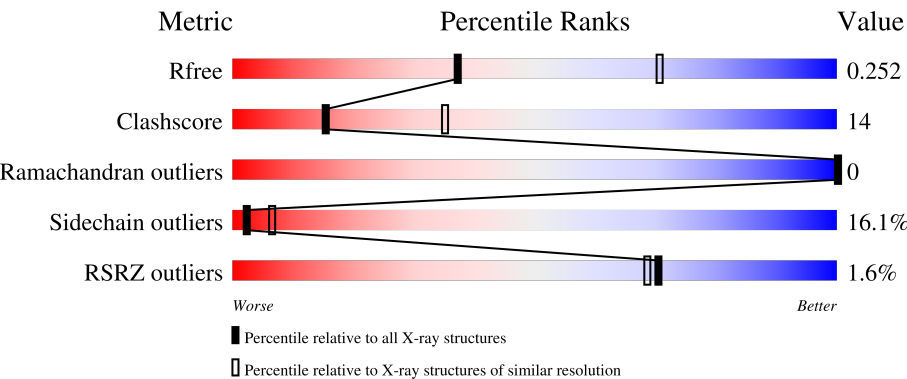
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry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

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.70 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	3538 (2.70-2.70)
Clashscore	190562	3843 (2.70-2.70)
Ramachandran outliers	187476	3778 (2.70-2.70)
Sidechain outliers	187428	3778 (2.70-2.70)
RSRZ outliers	180081	3538 (2.70-2.70)

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 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	B	115	2% 68% 29% .
1	E	115	% 64% 32% .
1	G	115	2% 70% 26% .
1	H	115	3% 60% 31% 8% .
2	A	222	36% 18% . 42%

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Mol	Chain	Length	Quality of chain
2	C	222	36%17%5%41%
2	D	222	36%20%•41%
2	F	222	2%37%17%5%41%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7570 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 NbSOS2 (14509).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	G	115	Total	C	N	O	S	0	0	0
			871	539	159	169	4			
1	E	115	Total	C	N	O	S	0	0	0
			871	539	159	169	4			
1	B	115	Total	C	N	O	S	0	0	0
			871	539	159	169	4			
1	H	114	Total	C	N	O	S	0	0	0
			867	537	158	168	4			

- Molecule 2 is a protein called LexA repressor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	A	129	Total	C	N	O	S	0	0	0
			1001	633	178	187	3			
2	C	130	Total	C	N	O	S	0	0	0
			1010	638	179	190	3			
2	D	130	Total	C	N	O	S	0	0	0
			1010	638	179	190	3			
2	F	131	Total	C	N	O	S	0	0	0
			1019	643	180	193	3			

There are 80 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP A0A1X3HXXW2
A	-18	GLY	-	expression tag	UNP A0A1X3HXXW2
A	-17	SER	-	expression tag	UNP A0A1X3HXXW2
A	-16	SER	-	expression tag	UNP A0A1X3HXXW2
A	-15	HIS	-	expression tag	UNP A0A1X3HXXW2
A	-14	HIS	-	expression tag	UNP A0A1X3HXXW2
A	-13	HIS	-	expression tag	UNP A0A1X3HXXW2
A	-12	HIS	-	expression tag	UNP A0A1X3HXXW2

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-11	HIS	-	expression tag	UNP A0A1X3HXLW2
A	-10	HIS	-	expression tag	UNP A0A1X3HXLW2
A	-9	SER	-	expression tag	UNP A0A1X3HXLW2
A	-8	SER	-	expression tag	UNP A0A1X3HXLW2
A	-7	GLY	-	expression tag	UNP A0A1X3HXLW2
A	-6	LEU	-	expression tag	UNP A0A1X3HXLW2
A	-5	VAL	-	expression tag	UNP A0A1X3HXLW2
A	-4	PRO	-	expression tag	UNP A0A1X3HXLW2
A	-3	ARG	-	expression tag	UNP A0A1X3HXLW2
A	-2	GLY	-	expression tag	UNP A0A1X3HXLW2
A	-1	SER	-	expression tag	UNP A0A1X3HXLW2
A	0	HIS	-	expression tag	UNP A0A1X3HXLW2
C	-19	MET	-	initiating methionine	UNP A0A1X3HXLW2
C	-18	GLY	-	expression tag	UNP A0A1X3HXLW2
C	-17	SER	-	expression tag	UNP A0A1X3HXLW2
C	-16	SER	-	expression tag	UNP A0A1X3HXLW2
C	-15	HIS	-	expression tag	UNP A0A1X3HXLW2
C	-14	HIS	-	expression tag	UNP A0A1X3HXLW2
C	-13	HIS	-	expression tag	UNP A0A1X3HXLW2
C	-12	HIS	-	expression tag	UNP A0A1X3HXLW2
C	-11	HIS	-	expression tag	UNP A0A1X3HXLW2
C	-10	HIS	-	expression tag	UNP A0A1X3HXLW2
C	-9	SER	-	expression tag	UNP A0A1X3HXLW2
C	-8	SER	-	expression tag	UNP A0A1X3HXLW2
C	-7	GLY	-	expression tag	UNP A0A1X3HXLW2
C	-6	LEU	-	expression tag	UNP A0A1X3HXLW2
C	-5	VAL	-	expression tag	UNP A0A1X3HXLW2
C	-4	PRO	-	expression tag	UNP A0A1X3HXLW2
C	-3	ARG	-	expression tag	UNP A0A1X3HXLW2
C	-2	GLY	-	expression tag	UNP A0A1X3HXLW2
C	-1	SER	-	expression tag	UNP A0A1X3HXLW2
C	0	HIS	-	expression tag	UNP A0A1X3HXLW2
D	-19	MET	-	initiating methionine	UNP A0A1X3HXLW2
D	-18	GLY	-	expression tag	UNP A0A1X3HXLW2
D	-17	SER	-	expression tag	UNP A0A1X3HXLW2
D	-16	SER	-	expression tag	UNP A0A1X3HXLW2
D	-15	HIS	-	expression tag	UNP A0A1X3HXLW2
D	-14	HIS	-	expression tag	UNP A0A1X3HXLW2
D	-13	HIS	-	expression tag	UNP A0A1X3HXLW2
D	-12	HIS	-	expression tag	UNP A0A1X3HXLW2
D	-11	HIS	-	expression tag	UNP A0A1X3HXLW2
D	-10	HIS	-	expression tag	UNP A0A1X3HXLW2

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Chain	Residue	Modelled	Actual	Comment	Reference
D	-9	SER	-	expression tag	UNP A0A1X3HXXW2
D	-8	SER	-	expression tag	UNP A0A1X3HXXW2
D	-7	GLY	-	expression tag	UNP A0A1X3HXXW2
D	-6	LEU	-	expression tag	UNP A0A1X3HXXW2
D	-5	VAL	-	expression tag	UNP A0A1X3HXXW2
D	-4	PRO	-	expression tag	UNP A0A1X3HXXW2
D	-3	ARG	-	expression tag	UNP A0A1X3HXXW2
D	-2	GLY	-	expression tag	UNP A0A1X3HXXW2
D	-1	SER	-	expression tag	UNP A0A1X3HXXW2
D	0	HIS	-	expression tag	UNP A0A1X3HXXW2
F	-19	MET	-	initiating methionine	UNP A0A1X3HXXW2
F	-18	GLY	-	expression tag	UNP A0A1X3HXXW2
F	-17	SER	-	expression tag	UNP A0A1X3HXXW2
F	-16	SER	-	expression tag	UNP A0A1X3HXXW2
F	-15	HIS	-	expression tag	UNP A0A1X3HXXW2
F	-14	HIS	-	expression tag	UNP A0A1X3HXXW2
F	-13	HIS	-	expression tag	UNP A0A1X3HXXW2
F	-12	HIS	-	expression tag	UNP A0A1X3HXXW2
F	-11	HIS	-	expression tag	UNP A0A1X3HXXW2
F	-10	HIS	-	expression tag	UNP A0A1X3HXXW2
F	-9	SER	-	expression tag	UNP A0A1X3HXXW2
F	-8	SER	-	expression tag	UNP A0A1X3HXXW2
F	-7	GLY	-	expression tag	UNP A0A1X3HXXW2
F	-6	LEU	-	expression tag	UNP A0A1X3HXXW2
F	-5	VAL	-	expression tag	UNP A0A1X3HXXW2
F	-4	PRO	-	expression tag	UNP A0A1X3HXXW2
F	-3	ARG	-	expression tag	UNP A0A1X3HXXW2
F	-2	GLY	-	expression tag	UNP A0A1X3HXXW2
F	-1	SER	-	expression tag	UNP A0A1X3HXXW2
F	0	HIS	-	expression tag	UNP A0A1X3HXXW2

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



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

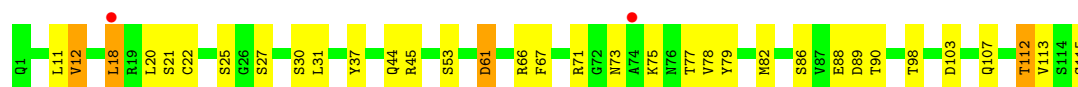
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	G	2	Total	O	0	0
			2	2		
4	A	5	Total	O	0	0
			5	5		
4	C	8	Total	O	0	0
			8	8		
4	E	12	Total	O	0	0
			12	12		
4	B	2	Total	O	0	0
			2	2		
4	D	7	Total	O	0	0
			7	7		
4	F	8	Total	O	0	0
			8	8		
4	H	2	Total	O	0	0
			2	2		

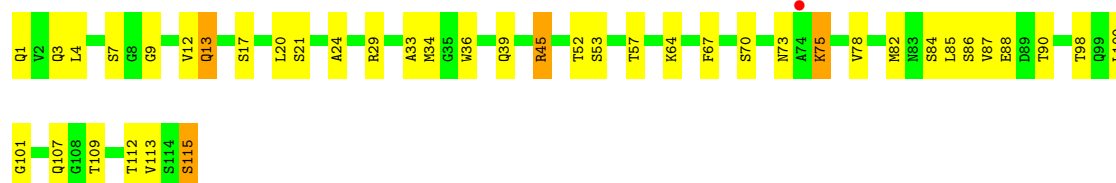
3 Residue-property plots [i](#)

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.

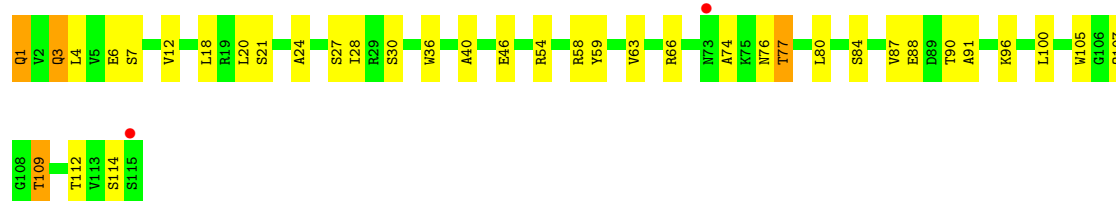
- Molecule 1: NbSOS2 (14509)



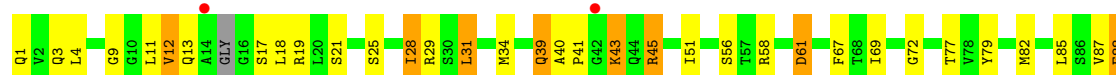
- Molecule 1: NbSOS2 (14509)



- Molecule 1: NbSOS2 (14509)



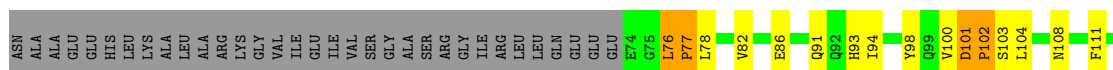
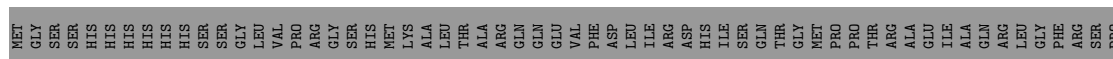
- Molecule 1: NbSOS2 (14509)





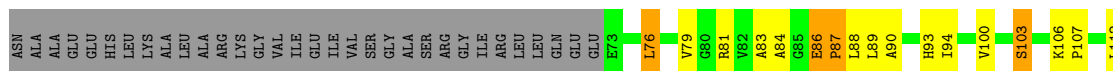
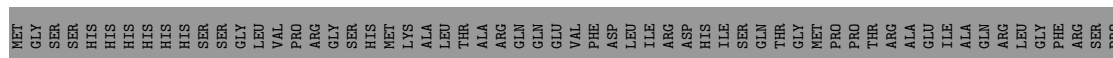
• Molecule 2: LexA repressor

Chain A: 36% 18% • 42%



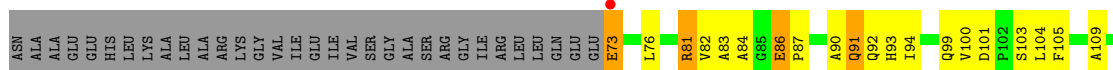
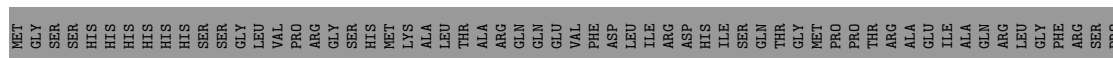
• Molecule 2: LexA repressor

Chain C: 36% 17% 5% 41%



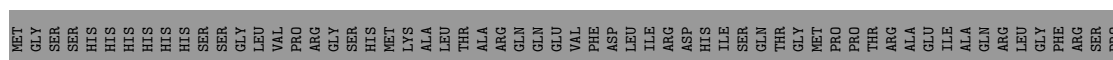
• Molecule 2: LexA repressor

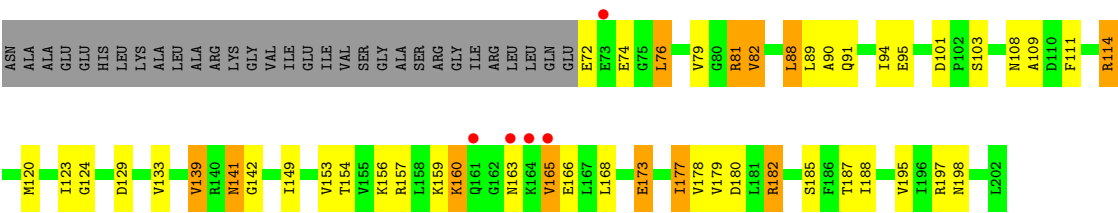
Chain D: 36% 20% • 41%



• Molecule 2: LexA repressor

Chain F: 2% 37% 17% 5% 41%





4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	112.67Å 112.67Å 144.75Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.25 – 2.70 48.25 – 2.70	Depositor EDS
% Data completeness (in resolution range)	96.1 (48.25-2.70) 96.1 (48.25-2.70)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.58 (at 2.69Å)	Xtriage
Refinement program	PHENIX 1.18.2	Depositor
R, R_{free}	0.228 , 0.260 0.225 , 0.252	Depositor DCC
R_{free} test set	1322 reflections (4.45%)	wwPDB-VP
Wilson B-factor (Å ²)	51.0	Xtriage
Anisotropy	0.031	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 12.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.037 for -h,-k,l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	7570	wwPDB-VP
Average B, all atoms (Å ²)	47.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.98% 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

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	B	0.57	0/884	0.80	0/1192
1	E	0.79	1/884 (0.1%)	1.05	0/1192
1	G	0.60	0/884	0.89	1/1192 (0.1%)
1	H	0.57	0/879	0.87	1/1184 (0.1%)
2	A	0.77	2/1016 (0.2%)	1.15	6/1373 (0.4%)
2	C	0.68	1/1025 (0.1%)	0.97	1/1385 (0.1%)
2	D	0.61	0/1025	0.92	0/1385
2	F	0.63	0/1034	0.96	0/1397
All	All	0.66	4/7631 (0.1%)	0.96	9/10300 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	77	PRO	C-O	-6.63	1.16	1.23
2	A	102	PRO	C-O	-5.96	1.16	1.24
2	C	87	PRO	C-O	-5.92	1.16	1.23
1	E	9	GLY	CA-C	-5.51	1.46	1.52

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	163	ASN	N-CA-C	-7.43	103.12	113.37
1	G	73	ASN	N-CA-C	-6.28	104.88	112.54
2	A	104	LEU	N-CA-C	-5.82	105.95	113.16
2	A	182	ARG	N-CA-C	-5.75	106.24	113.72
2	C	182	ARG	N-CA-C	-5.54	106.56	113.38
2	A	170	GLU	N-CA-C	-5.50	102.01	109.54
2	A	101	ASP	CA-CB-CG	5.42	118.02	112.60
1	H	89	ASP	N-CA-C	-5.08	107.37	113.21
2	A	77	PRO	N-CA-CB	-5.02	99.19	103.36

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	871	0	859	18	0
1	E	871	0	859	19	0
1	G	871	0	859	20	0
1	H	867	0	855	22	0
2	A	1001	0	1024	34	0
2	C	1010	0	1030	32	0
2	D	1010	0	1030	39	0
2	F	1019	0	1036	37	0
3	A	4	0	6	0	0
4	A	5	0	0	0	0
4	B	2	0	0	1	0
4	C	8	0	0	0	0
4	D	7	0	0	1	0
4	E	12	0	0	1	0
4	F	8	0	0	0	0
4	G	2	0	0	0	0
4	H	2	0	0	0	0
All	All	7570	0	7558	204	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 14.

All (204) 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:H:90:THR:HG23	1:H:112:THR:HA	1.61	0.81
2:C:141:ASN:HD21	2:C:160:LYS:H	1.29	0.81
2:F:160:LYS:HA	2:F:165:VAL:HG22	1.64	0.79
2:D:180:ASP:HB3	2:D:183:GLN:HE21	1.49	0.77
2:D:87:PRO:HG2	2:D:90:ALA:HB2	1.67	0.77
2:D:86:GLU:HB3	2:D:93:HIS:CE1	2.20	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:156:LYS:HE2	2:F:177:ILE:HD11	1.66	0.76
1:G:20:LEU:HG	1:G:82:MET:HE2	1.69	0.75
2:D:159:LYS:HB3	2:D:166:GLU:HB3	1.71	0.72
1:G:71:ARG:HG2	1:G:78:VAL:HG12	1.71	0.71
1:H:93:TYR:O	1:H:109:THR:HG22	1.90	0.70
1:H:9:GLY:HA2	1:H:18:LEU:HD13	1.73	0.70
2:C:138:ASP:OD1	2:C:140:ARG:NH2	2.25	0.69
2:C:149:ILE:HD11	2:C:167:LEU:HD12	1.74	0.69
2:C:129:ASP:OD1	2:C:197:ARG:HD3	1.92	0.69
2:F:120:MET:HE3	2:F:123:ILE:HD11	1.72	0.69
2:A:86:GLU:HB3	2:A:93:HIS:HE1	1.58	0.69
2:D:149:ILE:HD11	2:D:167:LEU:HD12	1.75	0.68
2:A:162:GLY:C	2:A:164:LYS:H	2.02	0.68
2:F:81:ARG:HH11	2:F:114:ARG:HE	1.46	0.64
1:G:18:LEU:C	1:G:18:LEU:HD22	2.23	0.63
1:H:69:ILE:HA	1:H:79:TYR:O	1.99	0.63
2:C:81:ARG:HD2	2:F:185:SER:HB2	1.80	0.63
2:C:84:ALA:HB3	2:F:163:ASN:HD21	1.64	0.62
1:E:34:MET:HB2	1:E:78:VAL:HG11	1.80	0.62
1:G:31:LEU:HD13	2:A:108:ASN:ND2	2.13	0.62
2:D:159:LYS:HE3	2:D:168:LEU:HD11	1.81	0.62
2:C:87:PRO:HG2	2:C:90:ALA:HB2	1.83	0.61
2:A:91:GLN:HA	2:A:94:ILE:HD12	1.83	0.61
2:A:197:ARG:NH2	2:C:124:GLY:O	2.34	0.60
1:G:18:LEU:CD1	1:G:82:MET:HE3	2.32	0.60
1:G:11:LEU:HD12	1:G:112:THR:HB	1.83	0.59
2:A:141:ASN:HB2	2:A:160:LYS:H	1.68	0.59
2:A:122:ASP:HB2	2:A:170:GLU:O	2.03	0.59
2:D:82:VAL:HG11	2:D:153:VAL:HG23	1.84	0.59
2:F:88:LEU:HD11	2:F:153:VAL:HG21	1.85	0.58
1:B:1:GLN:NE2	1:B:3:GLN:HE22	2.02	0.57
1:B:6:GLU:OE1	1:B:109:THR:HG22	2.05	0.57
2:F:142:GLY:O	2:F:157:ARG:NH1	2.37	0.57
2:D:197:ARG:HH22	2:F:124:GLY:C	2.13	0.56
1:H:17:SER:HA	1:H:82:MET:O	2.05	0.56
2:F:129:ASP:OD1	2:F:197:ARG:HD3	2.05	0.56
1:G:18:LEU:HD13	1:G:82:MET:HE3	1.88	0.56
2:C:83:ALA:HB3	2:C:86:GLU:HG3	1.88	0.56
1:B:54:ARG:HH22	2:D:73:GLU:HA	1.71	0.56
2:A:86:GLU:CB	2:A:93:HIS:HE1	2.19	0.56
2:D:86:GLU:HB3	2:D:93:HIS:HE1	1.66	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:200:ASP:HB3	4:E:202:HOH:O	2.05	0.55
2:F:129:ASP:OD2	2:F:197:ARG:NH1	2.39	0.55
2:D:180:ASP:HB3	2:D:183:GLN:NE2	2.19	0.55
2:C:112:LEU:HD23	2:C:132:ALA:HA	1.88	0.55
1:H:12:VAL:HG11	1:H:85:LEU:HD13	1.88	0.55
2:A:162:GLY:C	2:A:164:LYS:N	2.64	0.55
2:A:124:GLY:C	2:C:197:ARG:HH22	2.14	0.55
2:A:121:LYS:HE3	2:A:122:ASP:OD1	2.08	0.54
1:B:90:THR:HG23	1:B:112:THR:HA	1.90	0.54
1:E:12:VAL:O	1:E:113:VAL:HA	2.07	0.54
1:B:96:LYS:HD2	1:B:105:TRP:CH2	2.42	0.54
2:A:165:VAL:O	2:A:179:VAL:HG12	2.07	0.53
2:F:154:THR:HG1	2:F:156:LYS:HZ3	1.55	0.53
2:D:83:ALA:H	2:D:93:HIS:CE1	2.25	0.53
2:C:141:ASN:ND2	2:C:160:LYS:H	2.03	0.53
1:G:61:ASP:OD1	1:G:61:ASP:N	2.34	0.53
2:C:79:VAL:HG22	2:C:94:ILE:HD13	1.91	0.52
1:G:90:THR:HG23	1:G:112:THR:HA	1.92	0.52
1:E:29:ARG:HH21	1:E:100:LEU:HD21	1.75	0.52
2:F:141:ASN:HB2	2:F:160:LYS:HG2	1.91	0.52
2:A:86:GLU:HB3	2:A:93:HIS:CE1	2.43	0.51
2:F:159:LYS:HB2	2:F:168:LEU:HD11	1.93	0.51
2:A:181:LEU:CD2	2:A:186:PHE:HB3	2.39	0.51
1:B:107:GLN:HG2	4:B:201:HOH:O	2.09	0.51
1:E:67:PHE:CD1	1:E:82:MET:HA	2.46	0.50
2:D:129:ASP:OD1	2:D:197:ARG:HD3	2.11	0.50
2:C:103:SER:HB2	1:E:52:THR:HB	1.94	0.50
2:F:88:LEU:HD11	2:F:153:VAL:CG2	2.41	0.50
2:C:107:PRO:HG3	1:E:101:GLY:HA3	1.92	0.49
1:B:66:ARG:NH1	1:B:84:SER:O	2.45	0.49
2:D:135:LYS:HE2	2:D:135:LYS:C	2.37	0.49
2:F:82:VAL:HG11	2:F:88:LEU:HG	1.94	0.49
2:A:77:PRO:HD2	2:A:111:PHE:HA	1.95	0.49
1:B:54:ARG:HB2	2:D:101:ASP:OD2	2.13	0.49
2:D:200:ASP:OD1	1:H:58:ARG:NH1	2.46	0.49
2:C:120:MET:HE3	2:C:123:ILE:HD11	1.94	0.49
1:E:39:GLN:HB2	1:E:45:ARG:HD2	1.93	0.49
1:G:31:LEU:HD13	2:A:108:ASN:HD22	1.75	0.49
2:F:139:VAL:HG21	2:F:188:ILE:HG21	1.94	0.48
2:F:180:ASP:C	2:F:182:ARG:H	2.21	0.48
2:F:129:ASP:CG	2:F:197:ARG:HH11	2.22	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:67:PHE:CD1	1:H:82:MET:HA	2.49	0.48
1:G:67:PHE:CD1	1:G:82:MET:HA	2.49	0.48
1:E:75:LYS:HE3	1:E:75:LYS:HB2	1.54	0.48
1:G:66:ARG:NH1	1:G:89:ASP:OD2	2.45	0.48
2:F:76:LEU:HD21	2:F:109:ALA:HB1	1.96	0.48
2:F:163:ASN:ND2	2:F:182:ARG:HG2	2.29	0.48
1:H:4:LEU:HD12	1:H:105:TRP:C	2.39	0.47
2:A:146:VAL:HG12	2:A:189:GLU:HB2	1.96	0.47
2:A:179:VAL:HG21	2:A:186:PHE:CG	2.50	0.47
1:B:96:LYS:HD2	1:B:105:TRP:CZ2	2.48	0.47
1:E:36:TRP:CD1	1:E:36:TRP:N	2.83	0.47
1:B:4:LEU:HD23	1:B:24:ALA:HA	1.96	0.47
2:F:141:ASN:C	2:F:141:ASN:ND2	2.72	0.47
1:E:29:ARG:NH2	1:E:100:LEU:HD21	2.29	0.47
1:E:82:MET:HB3	1:E:85:LEU:HD21	1.96	0.47
1:E:86:SER:C	1:E:88:GLU:H	2.22	0.47
2:D:160:LYS:HG3	2:D:165:VAL:HG22	1.97	0.47
2:F:141:ASN:C	2:F:141:ASN:HD22	2.22	0.47
1:H:28:ILE:HD13	1:H:28:ILE:H	1.79	0.47
2:A:100:VAL:HG22	2:C:100:VAL:HG13	1.97	0.47
2:A:159:LYS:O	2:A:165:VAL:HA	2.14	0.47
2:A:180:ASP:C	2:A:182:ARG:H	2.21	0.47
2:C:168:LEU:CD2	2:C:176:PRO:HB3	2.45	0.47
2:D:103:SER:C	2:D:105:PHE:H	2.22	0.47
1:B:36:TRP:CZ3	1:B:80:LEU:HB2	2.50	0.47
2:F:90:ALA:O	2:F:91:GLN:C	2.58	0.47
1:E:87:VAL:O	1:E:90:THR:HG22	2.16	0.46
2:C:168:LEU:HD23	2:C:176:PRO:HB3	1.98	0.46
1:H:51:ILE:HA	1:H:56:SER:O	2.16	0.46
1:G:98:THR:HG22	1:G:103:ASP:HA	1.98	0.46
1:H:40:ALA:O	1:H:41:PRO:C	2.58	0.46
2:D:81:ARG:HA	2:D:81:ARG:HD3	1.33	0.46
2:F:141:ASN:OD1	2:F:160:LYS:N	2.45	0.46
2:C:165:VAL:CG2	2:C:179:VAL:HG13	2.46	0.46
2:F:89:LEU:HA	2:F:94:ILE:HD11	1.98	0.46
2:F:165:VAL:O	2:F:179:VAL:HG12	2.16	0.46
2:F:173:GLU:H	2:F:173:GLU:HG3	1.63	0.46
2:C:180:ASP:C	2:C:182:ARG:H	2.23	0.46
1:B:7:SER:OG	1:B:21:SER:OG	2.30	0.45
1:H:29:ARG:HD3	1:H:29:ARG:HA	1.63	0.45
2:D:76:LEU:HD11	2:D:109:ALA:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:24:ALA:HB3	1:B:76:ASN:HB3	1.98	0.45
2:A:165:VAL:HG12	2:A:179:VAL:CG1	2.46	0.45
2:C:146:VAL:HG22	2:C:155:VAL:HG22	1.99	0.45
1:B:74:ALA:HB3	1:B:77:THR:HB	1.99	0.45
2:D:122:ASP:HB2	2:D:170:GLU:O	2.17	0.45
1:H:40:ALA:HB3	1:H:43:LYS:HB2	1.98	0.45
2:C:165:VAL:HG23	2:C:179:VAL:HG13	1.99	0.44
2:F:81:ARG:NH1	2:F:114:ARG:HE	2.13	0.44
2:A:78:LEU:HD11	2:A:130:LEU:HD22	2.00	0.44
2:C:118:MET:O	2:C:121:LYS:HG2	2.17	0.44
1:B:40:ALA:HB2	1:B:91:ALA:HB2	1.98	0.44
1:E:13:GLN:OE1	1:E:115:SER:HA	2.18	0.44
1:E:34:MET:HE2	1:E:34:MET:HB3	1.74	0.44
2:D:139:VAL:HG12	2:D:191:LEU:HB2	2.00	0.44
2:A:76:LEU:HA	2:A:77:PRO:HD3	1.66	0.44
2:A:155:VAL:O	2:A:156:LYS:HG2	2.17	0.44
2:A:165:VAL:HG12	2:A:179:VAL:HG11	1.98	0.44
2:C:163:ASN:HB2	2:C:181:LEU:HB2	1.99	0.44
2:D:103:SER:C	2:D:105:PHE:N	2.76	0.44
1:B:59:TYR:HB3	1:B:63:VAL:HG21	2.00	0.44
2:F:120:MET:SD	2:F:156:LYS:HA	2.58	0.44
1:E:4:LEU:HD23	1:E:24:ALA:HA	1.99	0.43
2:D:149:ILE:HD11	2:D:167:LEU:CD1	2.45	0.43
2:F:141:ASN:OD1	2:F:159:LYS:HA	2.17	0.43
1:E:33:ALA:HB3	1:E:98:THR:HG22	1.99	0.43
2:D:76:LEU:HD12	2:D:76:LEU:HA	1.77	0.43
2:F:139:VAL:CG2	2:F:188:ILE:HG21	2.48	0.43
1:H:61:ASP:OD1	1:H:61:ASP:N	2.47	0.43
2:D:84:ALA:C	2:D:86:GLU:N	2.76	0.43
2:D:84:ALA:C	2:D:86:GLU:H	2.26	0.43
1:B:100:LEU:HD23	1:B:100:LEU:HA	1.82	0.43
2:C:136:THR:O	2:C:190:GLY:HA2	2.18	0.43
1:G:37:TYR:HB3	1:G:45:ARG:HG2	2.00	0.42
1:G:37:TYR:HE1	2:C:202:LEU:HG	1.84	0.42
1:B:58:ARG:HG3	1:B:59:TYR:N	2.33	0.42
2:A:131:LEU:HD21	2:A:192:ALA:HB2	2.01	0.42
2:C:180:ASP:C	2:C:182:ARG:N	2.77	0.42
2:C:86:GLU:HB2	2:C:93:HIS:CE1	2.54	0.42
1:H:88:GLU:C	1:H:90:THR:H	2.27	0.42
2:A:155:VAL:C	2:A:156:LYS:HG2	2.44	0.42
2:D:118:MET:HE2	2:D:127:ASP:HB3	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:98:TYR:HE2	2:A:198:ASN:HD21	1.61	0.42
2:D:91:GLN:C	2:D:93:HIS:H	2.28	0.42
2:D:100:VAL:HG11	2:D:196:ILE:HD11	2.02	0.42
2:A:160:LYS:HE2	2:A:160:LYS:HB2	1.48	0.42
1:H:34:MET:HG2	1:H:97:GLN:HG2	2.02	0.42
1:G:12:VAL:O	1:G:113:VAL:HA	2.20	0.42
2:D:104:LEU:O	2:F:198:ASN:HB2	2.19	0.42
2:D:180:ASP:CB	2:D:183:GLN:HE21	2.26	0.42
2:F:157:ARG:HB3	2:F:168:LEU:HB2	2.02	0.42
1:G:21:SER:HB2	1:G:79:TYR:HD2	1.85	0.41
2:D:82:VAL:HG11	2:D:153:VAL:CG2	2.49	0.41
1:H:39:GLN:NE2	1:H:45:ARG:HD2	2.36	0.41
1:H:72:GLY:HA3	1:H:79:TYR:HE1	1.85	0.41
2:D:83:ALA:HB3	2:D:86:GLU:HG2	2.03	0.41
2:D:101:ASP:HB3	2:D:104:LEU:CD1	2.51	0.41
2:A:101:ASP:HA	2:A:102:PRO:HD3	1.78	0.41
2:A:181:LEU:HD22	2:A:186:PHE:HB3	2.01	0.41
2:C:173:GLU:HG2	2:C:174:PHE:CD1	2.55	0.41
2:A:82:VAL:HG23	2:A:93:HIS:CE1	2.56	0.41
1:E:86:SER:C	1:E:88:GLU:N	2.79	0.41
2:F:180:ASP:C	2:F:182:ARG:N	2.79	0.41
1:G:21:SER:HB2	1:G:79:TYR:CD2	2.56	0.41
2:C:175:LYS:HD3	1:H:102:TYR:OH	2.20	0.41
2:D:99:GLN:HB2	2:F:101:ASP:HB2	2.02	0.41
2:D:119:SER:OG	2:D:156:LYS:NZ	2.43	0.41
1:H:4:LEU:HD13	1:H:95:CYS:SG	2.61	0.41
2:C:76:LEU:HD12	2:C:76:LEU:HA	1.82	0.41
1:H:31:LEU:HA	1:H:31:LEU:HD13	1.70	0.41
1:G:22:CYS:O	1:G:77:THR:HA	2.21	0.40
2:F:111:PHE:CZ	2:F:133:VAL:HG11	2.57	0.40
2:D:122:ASP:OD2	4:D:301:HOH:O	2.22	0.40
2:D:148:ARG:HD2	2:D:153:VAL:HG22	2.02	0.40
1:G:18:LEU:HD11	1:G:82:MET:HE3	2.02	0.40
1:E:33:ALA:HB3	1:E:98:THR:CG2	2.51	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	B	113/115 (98%)	100 (88%)	13 (12%)	0	100	100
1	E	113/115 (98%)	109 (96%)	4 (4%)	0	100	100
1	G	113/115 (98%)	108 (96%)	5 (4%)	0	100	100
1	H	110/115 (96%)	105 (96%)	5 (4%)	0	100	100
2	A	127/222 (57%)	117 (92%)	10 (8%)	0	100	100
2	C	128/222 (58%)	122 (95%)	6 (5%)	0	100	100
2	D	128/222 (58%)	115 (90%)	13 (10%)	0	100	100
2	F	129/222 (58%)	120 (93%)	9 (7%)	0	100	100
All	All	961/1348 (71%)	896 (93%)	65 (7%)	0	100	100

There are no Ramachandran outliers to report.

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	B	91/91 (100%)	77 (85%)	14 (15%)	2	7
1	E	91/91 (100%)	72 (79%)	19 (21%)	1	3
1	G	91/91 (100%)	77 (85%)	14 (15%)	2	7
1	H	91/91 (100%)	71 (78%)	20 (22%)	1	3
2	A	110/186 (59%)	97 (88%)	13 (12%)	5	13
2	C	111/186 (60%)	97 (87%)	14 (13%)	4	11

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	D	111/186 (60%)	98 (88%)	13 (12%)	5	13
2	F	112/186 (60%)	89 (80%)	23 (20%)	1	3
All	All	808/1108 (73%)	678 (84%)	130 (16%)	2	6

All (130) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	G	12	VAL
1	G	18	LEU
1	G	25	SER
1	G	27	SER
1	G	30	SER
1	G	44	GLN
1	G	53	SER
1	G	61	ASP
1	G	75	LYS
1	G	86	SER
1	G	88	GLU
1	G	107	GLN
1	G	112	THR
1	G	115	SER
2	A	76	LEU
2	A	103	SER
2	A	135	LYS
2	A	141	ASN
2	A	145	VAL
2	A	149	ILE
2	A	152	GLU
2	A	160	LYS
2	A	163	ASN
2	A	172	SER
2	A	178	VAL
2	A	185	SER
2	A	195	VAL
2	C	76	LEU
2	C	86	GLU
2	C	88	LEU
2	C	89	LEU
2	C	103	SER
2	C	106	LYS
2	C	130	LEU

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Mol	Chain	Res	Type
2	C	141	ASN
2	C	160	LYS
2	C	173	GLU
2	C	175	LYS
2	C	179	VAL
2	C	181	LEU
2	C	195	VAL
1	E	1	GLN
1	E	3	GLN
1	E	7	SER
1	E	13	GLN
1	E	17	SER
1	E	20	LEU
1	E	21	SER
1	E	45	ARG
1	E	53	SER
1	E	57	THR
1	E	64	LYS
1	E	70	SER
1	E	73	ASN
1	E	75	LYS
1	E	84	SER
1	E	107	GLN
1	E	109	THR
1	E	112	THR
1	E	115	SER
1	B	1	GLN
1	B	3	GLN
1	B	12	VAL
1	B	18	LEU
1	B	20	LEU
1	B	27	SER
1	B	28	ILE
1	B	30	SER
1	B	46	GLU
1	B	77	THR
1	B	87	VAL
1	B	88	GLU
1	B	109	THR
1	B	114	SER
2	D	73	GLU
2	D	81	ARG

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Mol	Chain	Res	Type
2	D	86	GLU
2	D	91	GLN
2	D	92	GLN
2	D	94	ILE
2	D	116	SER
2	D	121	LYS
2	D	135	LYS
2	D	158	LEU
2	D	172	SER
2	D	179	VAL
2	D	195	VAL
2	F	72	GLU
2	F	74	GLU
2	F	76	LEU
2	F	79	VAL
2	F	81	ARG
2	F	82	VAL
2	F	88	LEU
2	F	95	GLU
2	F	103	SER
2	F	108	ASN
2	F	114	ARG
2	F	139	VAL
2	F	141	ASN
2	F	149	ILE
2	F	160	LYS
2	F	165	VAL
2	F	166	GLU
2	F	173	GLU
2	F	177	ILE
2	F	178	VAL
2	F	182	ARG
2	F	187	THR
2	F	195	VAL
1	H	1	GLN
1	H	3	GLN
1	H	11	LEU
1	H	12	VAL
1	H	13	GLN
1	H	19	ARG
1	H	21	SER
1	H	25	SER

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Mol	Chain	Res	Type
1	H	28	ILE
1	H	31	LEU
1	H	39	GLN
1	H	43	LYS
1	H	45	ARG
1	H	61	ASP
1	H	77	THR
1	H	87	VAL
1	H	88	GLU
1	H	107	GLN
1	H	112	THR
1	H	114	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	G	44	GLN
1	G	76	ASN
1	G	107	GLN
2	A	92	GLN
2	A	99	GLN
2	A	161	GLN
2	A	183	GLN
2	C	97	HIS
2	C	99	GLN
2	C	108	ASN
2	C	141	ASN
2	C	143	GLN
2	C	161	GLN
2	C	183	GLN
1	B	1	GLN
1	B	13	GLN
1	B	83	ASN
1	B	110	GLN
2	D	91	GLN
2	D	93	HIS
2	D	99	GLN
2	D	183	GLN
2	F	91	GLN
2	F	99	GLN
2	F	163	ASN
1	H	39	GLN

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Mol	Chain	Res	Type
1	H	81	GLN
1	H	83	ASN

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](#)

1 ligand is 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	EDO	A	301	-	3,3,3	0.38	0	2,2,2	0.45	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	A	301	-	-	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	A	301	EDO	O1-C1-C2-O2

There are no ring outliers.

No monomer is involved in short contacts.

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	B	115/115 (100%)	0.22	2 (1%) 69 67	40, 49, 60, 70	0
1	E	115/115 (100%)	0.06	1 (0%) 81 80	37, 43, 52, 63	0
1	G	115/115 (100%)	0.19	2 (1%) 69 67	41, 51, 61, 66	0
1	H	114/115 (99%)	0.22	3 (2%) 57 54	44, 54, 67, 81	0
2	A	129/222 (58%)	0.03	1 (0%) 82 81	33, 40, 53, 65	0
2	C	130/222 (58%)	0.03	1 (0%) 82 81	33, 41, 55, 71	0
2	D	130/222 (58%)	0.25	1 (0%) 82 81	42, 51, 64, 66	0
2	F	131/222 (59%)	0.09	5 (3%) 44 40	39, 45, 57, 67	0
All	All	979/1348 (72%)	0.13	16 (1%) 70 68	33, 47, 62, 81	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	F	163	ASN	3.6
1	G	18	LEU	3.2
1	E	74	ALA	2.9
1	B	73	ASN	2.8
1	B	115	SER	2.7
1	H	115	SER	2.6
1	H	14	ALA	2.5
2	F	161	GLN	2.5
2	F	73	GLU	2.5
2	F	165	VAL	2.3
2	C	161	GLN	2.2
1	H	42	GLY	2.2
1	G	74	ALA	2.1
2	D	73	GLU	2.1
2	F	164	LYS	2.1
2	A	163	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 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	A	301	4/4	0.77	0.20	43,45,45,48	0

6.5 Other polymers [i](#)

There are no such residues in this entry.