



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 06:25 AM GMT

PDB ID : 2B7X  
Title : Sequential reorganization of beta-sheet topology by insertion of a single strand  
Authors : Sagermann, M.; Matthews, B.W.  
Deposited on : 2005-10-05  
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](mailto:validation@mail.wwpdb.org)  
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

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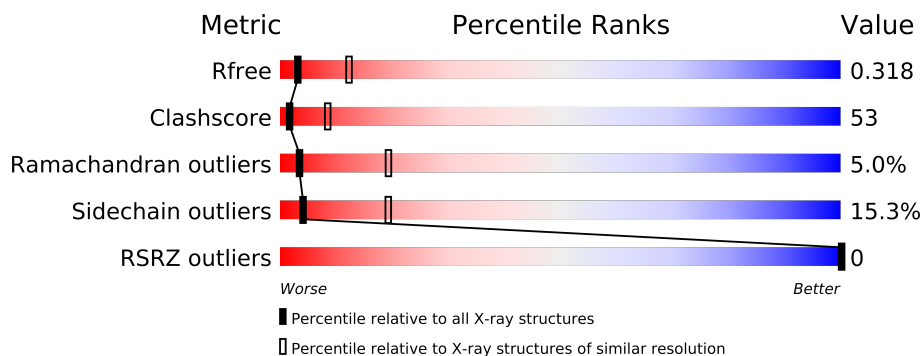
The following versions of software and data (see [references](#)) were used in the production of this report:

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  $\geq 3$ , 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	170	
1	B	170	
1	C	170	
1	D	170	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 5085 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 Lysozyme.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	157	Total	C	N	O	S	0	0	0
			1256	791	227	233	5			
1	B	157	Total	C	N	O	S	0	0	0
			1256	791	227	233	5			
1	C	157	Total	C	N	O	S	0	0	0
			1256	791	227	233	5			
1	D	160	Total	C	N	O	S	0	0	0
			1277	805	230	237	5			

There are 32 discrepancies between the modelled and reference sequences:

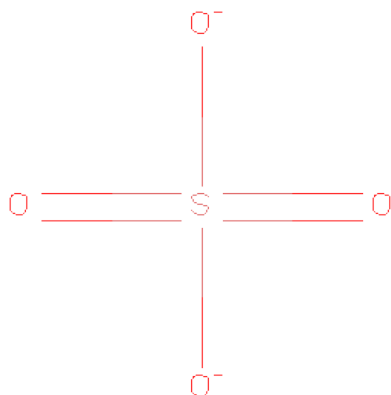
Chain	Residue	Modelled	Actual	Comment	Reference
A	31	TYR	-	INSERTION	UNP P00720
A	32	THR	-	INSERTION	UNP P00720
A	33	ILE	-	INSERTION	UNP P00720
A	34	GLY	-	INSERTION	UNP P00720
A	35	ILE	-	INSERTION	UNP P00720
A	36	GLY	-	INSERTION	UNP P00720
A	60	THR	CYS	ENGINEERED	UNP P00720
A	103	ALA	CYS	ENGINEERED	UNP P00720
B	25	TYR	-	INSERTION	UNP P00720
B	26	THR	-	INSERTION	UNP P00720
B	27	ILE	-	INSERTION	UNP P00720
B	28	GLY	-	INSERTION	UNP P00720
B	29	ILE	-	INSERTION	UNP P00720
B	30	GLY	-	INSERTION	UNP P00720
B	60	THR	CYS	ENGINEERED	UNP P00720
B	103	ALA	CYS	ENGINEERED	UNP P00720
C	31	TYR	-	INSERTION	UNP P00720
C	32	THR	-	INSERTION	UNP P00720
C	33	ILE	-	INSERTION	UNP P00720
C	34	GLY	-	INSERTION	UNP P00720
C	35	ILE	-	INSERTION	UNP P00720

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Chain	Residue	Modelled	Actual	Comment	Reference
C	36	GLY	-	INSERTION	UNP P00720
C	60	THR	CYS	ENGINEERED	UNP P00720
C	103	ALA	CYS	ENGINEERED	UNP P00720
D	31	TYR	-	INSERTION	UNP P00720
D	32	THR	-	INSERTION	UNP P00720
D	33	ILE	-	INSERTION	UNP P00720
D	34	GLY	-	INSERTION	UNP P00720
D	35	ILE	-	INSERTION	UNP P00720
D	36	GLY	-	INSERTION	UNP P00720
D	60	THR	CYS	ENGINEERED	UNP P00720
D	103	ALA	CYS	ENGINEERED	UNP P00720

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O<sub>4</sub>S).



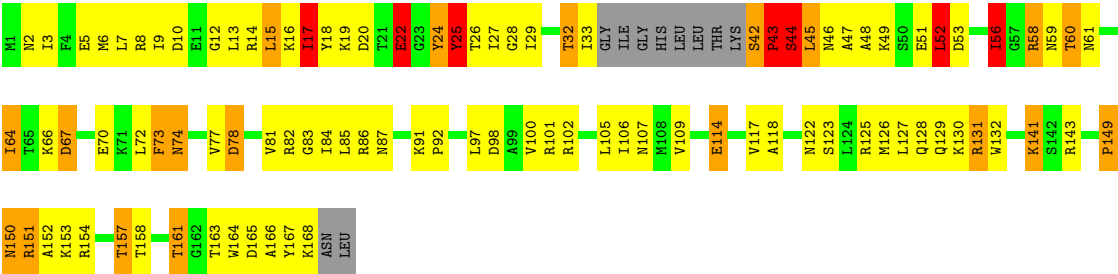
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	D	1	Total O S 5 4 1	0	0
2	D	1	Total O S 5 4 1	0	0
2	D	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0
2	C	1	Total O S 5 4 1	0	0

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		





## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	61.38Å 78.08Å 143.35Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 3.00 20.01 – 2.40	Depositor EDS
% Data completeness (in resolution range)	(Not available) (20.00-3.00) 70.1 (20.01-2.40)	Depositor EDS
$R_{merge}$	0.07	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.09 (at 2.41Å)	Xtriage
Refinement program	CNS	Depositor
R, $R_{free}$	0.239 , 0.319 0.237 , 0.318	Depositor DCC
$R_{free}$ test set	521 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	29.6	Xtriage
Anisotropy	0.188	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 4.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 19371 reflections (0.005%)	Xtriage
$F_o, F_c$ correlation	0.87	EDS
Total number of atoms	5085	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	8.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 38.10 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.8790e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

<sup>1</sup>Intensities estimated from amplitudes.



## 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: SO4

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	1.87	22/1274 (1.7%)	2.60	42/1714 (2.5%)
1	B	2.51	18/1274 (1.4%)	2.61	32/1714 (1.9%)
1	C	1.26	5/1274 (0.4%)	1.20	7/1714 (0.4%)
1	D	1.95	10/1296 (0.8%)	1.43	14/1745 (0.8%)
All	All	1.95	55/5118 (1.1%)	2.06	95/6887 (1.4%)

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	A	1	2
1	B	0	6
1	D	1	1
All	All	2	9

All (55) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	21	THR	CB-OG1	64.69	2.72	1.43
1	D	43	PRO	N-CD	31.03	1.91	1.47
1	D	43	PRO	CB-CG	27.65	2.88	1.50
1	D	44	SER	CB-OG	-25.73	1.08	1.42
1	A	60	THR	CB-OG1	-23.03	0.97	1.43
1	B	22	GLU	CB-CG	17.71	1.85	1.52
1	A	147	GLN	CB-CG	-17.56	1.05	1.52
1	A	150	ASN	CB-CG	-17.30	1.11	1.51
1	B	22	GLU	C-N	-17.16	1.02	1.33
1	B	131	ARG	C-N	-16.96	0.95	1.34
1	D	24	TYR	CA-CB	16.35	1.90	1.53

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	149	PRO	N-CD	15.98	1.70	1.47
1	C	100	VAL	CB-CG1	15.13	1.84	1.52
1	A	107	ASN	CB-CG	-14.82	1.17	1.51
1	B	19	LYS	CB-CG	14.22	1.91	1.52
1	D	25	TYR	CB-CG	-13.60	1.31	1.51
1	B	17	ILE	C-N	-13.51	1.02	1.34
1	B	132	TRP	C-N	-13.36	1.03	1.34
1	B	20	ASP	C-N	-12.84	1.04	1.34
1	A	106	ILE	CB-CG2	12.66	1.92	1.52
1	A	146	ASN	CB-CG	-10.72	1.26	1.51
1	D	58	ARG	CB-CG	-10.18	1.25	1.52
1	A	148	THR	CB-OG1	-9.88	1.23	1.43
1	A	60	THR	CB-CG2	9.22	1.82	1.52
1	B	131	ARG	CB-CG	-8.98	1.28	1.52
1	D	17	ILE	CB-CG2	8.52	1.79	1.52
1	A	26	THR	CB-OG1	8.41	1.60	1.43
1	A	21	THR	CB-OG1	-8.33	1.26	1.43
1	A	25	TYR	C-N	-8.03	1.15	1.34
1	C	100	VAL	CB-CG2	-8.00	1.36	1.52
1	B	130	LYS	CA-CB	7.99	1.71	1.53
1	B	128	GLN	C-N	-7.92	1.15	1.34
1	A	104	ALA	C-N	7.43	1.51	1.34
1	B	21	THR	C-N	-7.38	1.17	1.34
1	B	20	ASP	CB-CG	-7.00	1.37	1.51
1	A	18	TYR	CE2-CZ	6.99	1.47	1.38
1	D	58	ARG	C-N	6.97	1.50	1.34
1	A	146	ASN	C-N	-6.96	1.18	1.34
1	A	106	ILE	CB-CG1	-6.92	1.34	1.54
1	D	15	LEU	C-N	-6.32	1.19	1.34
1	A	20	ASP	C-N	6.24	1.48	1.34
1	D	42	SER	C-N	-6.19	1.22	1.34
1	B	18	TYR	CD1-CE1	6.07	1.48	1.39
1	B	21	THR	CB-CG2	-5.97	1.32	1.52
1	A	61	ASN	CB-CG	-5.79	1.37	1.51
1	A	149	PRO	CB-CG	-5.53	1.22	1.50
1	B	134	GLU	CG-CD	5.48	1.60	1.51
1	C	100	VAL	C-N	-5.47	1.21	1.34
1	A	102	ARG	CB-CG	-5.36	1.38	1.52
1	A	26	THR	CB-CG2	-5.34	1.34	1.52
1	B	70	GLU	CB-CG	-5.31	1.42	1.52
1	C	24	TYR	CE1-CZ	5.21	1.45	1.38
1	C	99	ALA	C-N	-5.20	1.22	1.34

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	24	TYR	CD2-CE2	5.04	1.47	1.39
1	B	18	TYR	CG-CD1	5.04	1.45	1.39

All (95) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	25	TYR	CB-CG-CD1	47.27	149.36	121.00
1	A	25	TYR	CB-CG-CD2	-46.86	92.88	121.00
1	B	20	ASP	CB-CG-OD1	32.83	147.85	118.30
1	A	24	TYR	CB-CG-CD2	-32.43	101.54	121.00
1	B	20	ASP	CB-CG-OD2	-30.70	90.67	118.30
1	B	132	TRP	CB-CG-CD2	-30.23	87.31	126.60
1	B	132	TRP	CB-CG-CD1	29.00	164.70	127.00
1	A	147	GLN	CA-CB-CG	24.93	168.25	113.40
1	B	22	GLU	O-C-N	-24.18	82.09	123.20
1	B	21	THR	CA-CB-OG1	-20.17	66.64	109.00
1	B	18	TYR	CB-CG-CD2	19.85	132.91	121.00
1	B	21	THR	OG1-CB-CG2	-19.80	64.45	110.00
1	B	21	THR	O-C-N	-19.71	91.16	122.70
1	B	22	GLU	C-N-CA	19.71	163.69	122.30
1	B	22	GLU	CA-C-N	19.39	154.97	116.20
1	B	132	TRP	O-C-N	-18.45	93.18	122.70
1	B	131	ARG	C-N-CA	17.71	165.99	121.70
1	A	24	TYR	CB-CG-CD1	17.66	131.59	121.00
1	A	146	ASN	CB-CG-OD1	-17.53	86.55	121.60
1	B	132	TRP	C-N-CA	16.88	163.89	121.70
1	B	18	TYR	CB-CG-CD1	-16.62	111.03	121.00
1	A	21	THR	CA-CB-CG2	-15.83	90.23	112.40
1	D	43	PRO	CA-CB-CG	-14.94	75.62	104.00
1	B	22	GLU	CB-CG-CD	-14.73	74.44	114.20
1	A	149	PRO	CA-N-CD	-14.47	91.24	111.50
1	A	102	ARG	CA-CB-CG	14.34	144.95	113.40
1	A	146	ASN	CB-CG-ND2	14.25	150.90	116.70
1	A	150	ASN	CA-CB-CG	13.97	144.13	113.40
1	B	20	ASP	CA-CB-CG	13.73	143.61	113.40
1	D	16	LYS	CB-CG-CD	13.53	146.78	111.60
1	D	43	PRO	N-CA-CB	13.15	119.08	103.30
1	B	21	THR	CA-C-N	12.93	145.64	117.20
1	B	131	ARG	CA-CB-CG	12.84	141.64	113.40
1	B	21	THR	C-N-CA	11.96	151.60	121.70
1	B	132	TRP	CA-C-N	11.90	143.38	117.20
1	A	148	THR	CA-CB-CG2	-11.42	96.41	112.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	146	ASN	CA-CB-CG	11.19	138.02	113.40
1	A	106	ILE	CA-CB-CG2	-10.72	89.45	110.90
1	C	97	LEU	CB-CG-CD1	-10.70	92.81	111.00
1	D	56	ILE	CB-CG1-CD1	-10.63	84.15	113.90
1	C	97	LEU	CB-CG-CD2	10.39	128.67	111.00
1	A	60	THR	CA-CB-OG1	10.37	130.77	109.00
1	D	25	TYR	CA-CB-CG	10.19	132.76	113.40
1	A	107	ASN	CA-CB-CG	10.05	135.52	113.40
1	B	131	ARG	O-C-N	-9.78	107.05	122.70
1	D	45	LEU	CB-CG-CD2	-9.77	94.39	111.00
1	D	42	SER	O-C-N	-9.66	102.75	121.10
1	A	22	GLU	CA-CB-CG	8.94	133.07	113.40
1	C	100	VAL	CA-CB-CG1	-8.92	97.53	110.90
1	C	100	VAL	CA-CB-CG2	8.71	123.97	110.90
1	A	107	ASN	CB-CG-ND2	-8.25	96.91	116.70
1	A	150	ASN	CB-CG-OD1	-8.24	105.12	121.60
1	B	53	ASP	CB-CG-OD2	8.12	125.61	118.30
1	D	56	ILE	CA-CB-CG1	8.10	126.38	111.00
1	A	149	PRO	CA-CB-CG	7.79	119.60	104.80
1	A	60	THR	CA-CB-CG2	-7.62	101.73	112.40
1	D	25	TYR	CB-CG-CD2	-7.35	116.59	121.00
1	D	17	ILE	CA-CB-CG1	7.33	124.92	111.00
1	A	21	THR	OG1-CB-CG2	7.03	126.16	110.00
1	B	21	THR	CA-CB-CG2	7.03	122.24	112.40
1	D	52	LEU	CB-CG-CD1	6.99	122.88	111.00
1	A	23	GLY	O-C-N	6.82	133.62	122.70
1	A	107	ASN	CB-CG-OD1	6.78	135.17	121.60
1	A	26	THR	CA-CB-CG2	6.76	121.86	112.40
1	A	102	ARG	CB-CG-CD	6.70	129.02	111.60
1	A	106	ILE	CG1-CB-CG2	-6.69	96.67	111.40
1	D	42	SER	CA-C-N	6.59	135.56	117.10
1	D	16	LYS	CA-CB-CG	6.59	127.89	113.40
1	B	132	TRP	N-CA-CB	6.42	122.16	110.60
1	B	19	LYS	CA-CB-CG	-6.40	99.31	113.40
1	A	26	THR	CA-CB-OG1	-6.35	95.67	109.00
1	A	60	THR	OG1-CB-CG2	6.22	124.32	110.00
1	A	150	ASN	CB-CG-ND2	6.09	131.32	116.70
1	A	21	THR	CA-CB-OG1	-6.08	96.23	109.00
1	B	131	ARG	CB-CG-CD	-6.04	95.90	111.60
1	C	86	ARG	NE-CZ-NH1	6.02	123.31	120.30
1	A	82	ARG	NE-CZ-NH2	-5.93	117.33	120.30
1	D	22	GLU	C-N-CA	5.92	134.73	122.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	24	TYR	CA-CB-CG	5.88	124.57	113.40
1	C	53	ASP	CB-CG-OD2	5.86	123.58	118.30
1	A	107	ASN	O-C-N	5.84	132.04	122.70
1	A	105	LEU	CB-CG-CD2	-5.84	101.08	111.00
1	A	61	ASN	CB-CG-OD1	-5.76	110.08	121.60
1	B	132	TRP	CB-CA-C	5.69	121.78	110.40
1	B	129	GLN	N-CA-CB	-5.67	100.40	110.60
1	A	25	TYR	CA-C-N	5.54	129.39	117.20
1	A	148	THR	OG1-CB-CG2	5.54	122.75	110.00
1	C	112	MET	CG-SD-CE	5.53	109.06	100.20
1	B	131	ARG	CA-C-N	5.52	129.35	117.20
1	B	129	GLN	CB-CA-C	5.51	121.41	110.40
1	B	22	GLU	CA-CB-CG	-5.47	101.38	113.40
1	A	59	ASN	O-C-N	-5.35	114.14	122.70
1	A	146	ASN	O-C-N	-5.23	114.34	122.70
1	A	106	ILE	CB-CG1-CD1	-5.16	99.46	113.90
1	A	107	ASN	C-N-CA	-5.03	109.13	121.70

All (2) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	60	THR	CB
1	D	56	ILE	CB

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	24	TYR	Sidechain
1	A	25	TYR	Mainchain
1	B	132	TRP	Peptide
1	B	20	ASP	Mainchain
1	B	21	THR	Mainchain,Peptide
1	B	22	GLU	Mainchain,Peptide
1	D	25	TYR	Sidechain

## 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	1256	0	1271	125	0
1	B	1256	0	1266	124	0
1	C	1256	0	1271	139	0
1	D	1277	0	1294	152	0
2	A	10	0	0	3	0
2	C	15	0	0	3	0
2	D	15	0	0	4	0
All	All	5085	0	5102	540	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 53.

All (540) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:17:ILE:CG2	1:D:17:ILE:CB	1.79	1.57
1:A:60:THR:CB	1:A:60:THR:CG2	1.82	1.53
1:C:100:VAL:CG1	1:C:100:VAL:CB	1.84	1.52
1:B:22:GLU:CG	1:B:22:GLU:CB	1.85	1.50
1:D:24:TYR:CB	1:D:24:TYR:CA	1.89	1.49
1:B:19:LYS:CG	1:B:19:LYS:CB	1.90	1.49
1:A:106:ILE:CG2	1:A:106:ILE:CB	1.92	1.48
1:A:149:PRO:N	1:A:149:PRO:CD	1.70	1.43
1:D:26:THR:O	1:D:27:ILE:CD1	1.72	1.38
1:D:43:PRO:CD	1:D:43:PRO:N	1.91	1.34
1:B:22:GLU:CD	1:B:22:GLU:CB	2.06	1.24
1:D:26:THR:C	1:D:27:ILE:HD12	1.59	1.21
1:D:26:THR:O	1:D:27:ILE:HD12	1.04	1.20
1:D:45:LEU:H	1:D:45:LEU:HD23	1.12	1.12
1:B:131:ARG:HE	1:B:131:ARG:HA	1.03	1.11
1:C:71:LYS:HE2	1:C:75:GLN:HE21	1.19	1.04
1:D:87:ASN:HB2	1:D:114:GLU:OE2	1.61	1.01
1:D:25:TYR:CE2	1:D:45:LEU:HB3	1.96	1.00
1:D:53:ASP:O	1:D:56:ILE:O	1.80	0.99
1:C:27:ILE:HD11	1:C:64:ILE:HD11	1.45	0.98
1:C:14:ARG:HE	1:C:19:LYS:HB2	1.28	0.98
1:D:32:THR:O	1:D:33:ILE:HG13	1.64	0.97
1:A:98:ASP:O	1:A:102:ARG:HG2	1.65	0.96
1:D:52:LEU:HG	1:D:60:THR:HG21	1.48	0.94
1:A:106:ILE:CG2	1:A:106:ILE:CA	2.45	0.93
1:A:106:ILE:CG1	1:A:106:ILE:CG2	2.46	0.93

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:142:SER:HB2	1:B:145:TYR:H	1.33	0.91
1:B:131:ARG:HA	1:B:131:ARG:NE	1.86	0.90
1:A:151:ARG:O	1:A:155:VAL:HG12	1.70	0.90
1:D:45:LEU:N	1:D:45:LEU:HD23	1.85	0.89
1:D:43:PRO:HG2	1:D:43:PRO:O	1.71	0.89
1:C:153:LYS:O	1:C:157:THR:HG23	1.71	0.89
1:B:22:GLU:CD	1:B:22:GLU:HB2	1.90	0.88
1:D:33:ILE:HG23	1:D:51:GLU:HG2	1.56	0.88
1:C:52:LEU:HG	1:C:60:THR:HG21	1.53	0.88
1:A:106:ILE:N	1:A:106:ILE:HD12	1.87	0.88
1:D:12:GLY:O	1:D:29:ILE:HA	1.74	0.87
1:D:13:LEU:HD22	1:D:14:ARG:H	1.39	0.87
1:B:21:THR:HG23	1:B:21:THR:OG1	1.74	0.86
1:B:22:GLU:OE1	1:B:22:GLU:CB	2.23	0.86
1:D:43:PRO:CG	1:D:43:PRO:O	2.24	0.86
1:C:1:MET:HG3	1:C:5:GLU:HB2	1.58	0.86
1:C:100:VAL:CG1	1:C:100:VAL:CA	2.55	0.85
1:B:137:VAL:HG12	1:B:141:LYS:NZ	1.91	0.85
1:A:106:ILE:H	1:A:106:ILE:HD12	1.40	0.84
1:D:17:ILE:CG1	1:D:17:ILE:CG2	2.54	0.84
1:C:25:TYR:O	1:C:32:THR:HG23	1.75	0.84
1:C:100:VAL:CG1	1:C:100:VAL:CG2	2.56	0.83
1:B:157:THR:O	1:B:161:THR:HG23	1.80	0.81
1:B:151:ARG:CA	1:B:154:ARG:HH21	1.94	0.81
1:A:138:ASN:OD1	1:A:141:LYS:NZ	2.14	0.80
1:C:10:ASP:OD2	1:C:151:ARG:HG3	1.80	0.80
1:C:86:ARG:HB3	1:C:86:ARG:NH1	1.96	0.80
1:C:18:TYR:HE1	1:C:20:ASP:HA	1.45	0.80
1:C:14:ARG:NE	1:C:19:LYS:HB2	1.95	0.80
1:A:60:THR:CA	1:A:60:THR:CG2	2.60	0.80
1:B:19:LYS:CG	1:B:19:LYS:CA	2.60	0.79
1:D:25:TYR:CZ	1:D:45:LEU:HB3	2.18	0.79
1:D:17:ILE:CA	1:D:17:ILE:CG2	2.61	0.79
1:B:22:GLU:OE1	1:B:22:GLU:HB2	1.82	0.79
1:D:157:THR:CG2	1:D:166:ALA:HB2	2.13	0.79
1:B:137:VAL:HG12	1:B:141:LYS:HZ2	1.48	0.78
1:B:151:ARG:HA	1:B:154:ARG:HH21	1.48	0.78
1:D:151:ARG:HH11	1:D:151:ARG:HG3	1.46	0.78
1:B:10:ASP:HB3	1:B:151:ARG:HD2	1.65	0.78
1:A:145:TYR:OH	1:A:153:LYS:HE3	1.84	0.78
1:C:142:SER:OG	1:C:144:TRP:HB3	1.84	0.77
1:C:27:ILE:CD1	1:C:64:ILE:HD11	2.15	0.77

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:48:ALA:O	1:D:51:GLU:HG3	1.85	0.76
1:C:52:LEU:HD12	1:C:56:ILE:HG22	1.66	0.76
1:A:111:GLN:HB2	1:A:151:ARG:NH2	2.01	0.75
1:B:139:LEU:O	1:B:142:SER:OG	2.04	0.75
1:B:93:VAL:HA	1:B:128:GLN:OE1	1.86	0.75
1:C:56:ILE:HD13	1:C:68:GLU:CD	2.06	0.74
1:C:72:LEU:O	1:C:73:PHE:C	2.22	0.74
1:D:17:ILE:CG2	1:D:18:TYR:N	2.50	0.74
1:B:22:GLU:CG	1:B:22:GLU:CA	2.66	0.74
1:D:26:THR:O	1:D:27:ILE:CG1	2.36	0.74
1:A:150:ASN:HB2	2:A:807:SO4:O4	1.87	0.74
1:D:15:LEU:O	1:D:64:ILE:HD13	1.86	0.73
1:B:107:ASN:HD21	1:B:151:ARG:HE	1.37	0.73
1:D:8:ARG:NH1	1:D:12:GLY:HA2	2.03	0.73
1:D:64:ILE:HD13	1:D:64:ILE:H	1.55	0.72
1:B:56:ILE:HD11	1:B:68:GLU:HB3	1.70	0.72
1:B:19:LYS:HG2	1:B:19:LYS:H	1.54	0.72
1:C:117:VAL:O	1:C:117:VAL:HG12	1.90	0.71
1:B:107:ASN:ND2	1:B:155:VAL:HG21	2.05	0.71
1:C:93:VAL:HG21	1:C:124:LEU:CD2	2.19	0.71
1:B:98:ASP:O	1:B:102:ARG:HB2	1.89	0.71
1:D:157:THR:HG21	1:D:166:ALA:HB2	1.72	0.71
1:D:74:ASN:O	1:D:77:VAL:HG12	1.89	0.71
1:C:53:ASP:OD1	1:C:59:ASN:HA	1.91	0.71
1:D:91:LYS:HB3	1:D:92:PRO:HD3	1.73	0.70
1:C:5:GLU:O	1:C:9:ILE:HD13	1.90	0.70
1:A:150:ASN:HB2	2:A:807:SO4:S	2.31	0.70
1:C:18:TYR:CE1	1:C:20:ASP:HA	2.27	0.70
1:D:153:LYS:O	1:D:157:THR:HB	1.92	0.70
1:A:29:ILE:O	1:A:30:GLY:C	2.30	0.70
1:B:22:GLU:C	1:B:22:GLU:CG	2.60	0.69
1:D:53:ASP:OD1	1:D:59:ASN:HA	1.90	0.69
1:D:13:LEU:HD22	1:D:14:ARG:N	2.08	0.69
1:A:103:ALA:HA	1:A:106:ILE:HD13	1.74	0.69
1:D:165:ASP:HA	1:D:168:LYS:HE2	1.73	0.69
1:D:161:THR:OG1	1:D:163:THR:OG1	2.09	0.69
1:B:93:VAL:HG21	1:B:124:LEU:HD22	1.73	0.69
1:A:102:ARG:O	1:A:106:ILE:HD12	1.93	0.69
1:A:76:ASP:O	1:A:79:ALA:HB3	1.93	0.68
1:A:104:ALA:HB1	1:A:155:VAL:HG23	1.73	0.68
1:C:143:ARG:NH1	1:C:147:GLN:HB2	2.08	0.68
1:B:79:ALA:HA	1:B:82:ARG:NH1	2.08	0.68

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:76:ASP:O	1:B:79:ALA:HB3	1.93	0.68
1:D:167:TYR:O	1:D:168:LYS:C	2.32	0.67
1:D:109:VAL:HA	1:D:117:VAL:HG21	1.75	0.67
1:A:3:ILE:O	1:A:6:MET:HB3	1.94	0.67
1:C:1:MET:CG	1:C:5:GLU:HB2	2.23	0.67
1:C:120:PHE:O	1:C:124:LEU:HB2	1.95	0.67
1:D:149:PRO:O	1:D:152:ALA:N	2.27	0.67
1:A:102:ARG:O	1:A:106:ILE:CD1	2.41	0.67
1:D:42:SER:C	1:D:43:PRO:CD	2.63	0.67
1:B:22:GLU:C	1:B:22:GLU:HG3	2.14	0.67
1:B:79:ALA:HA	1:B:82:ARG:HH12	1.60	0.67
1:B:84:ILE:HD13	1:B:90:LEU:HD13	1.76	0.67
1:C:29:ILE:HD12	1:C:29:ILE:N	2.10	0.66
1:D:122:ASN:HB2	2:D:802:SO4:O2	1.94	0.66
1:C:13:LEU:HD11	1:C:15:LEU:CD1	2.25	0.66
1:D:25:TYR:CE2	1:D:45:LEU:CB	2.75	0.66
1:B:22:GLU:O	1:B:22:GLU:HG3	1.95	0.66
1:B:104:ALA:HB3	1:B:159:PHE:CE1	2.31	0.66
1:A:60:THR:N	1:A:60:THR:CG2	2.58	0.66
1:A:98:ASP:OD2	1:A:98:ASP:C	2.35	0.66
1:C:86:ARG:CZ	1:C:86:ARG:HB3	2.26	0.66
1:C:150:ASN:HB2	2:C:805:SO4:O4	1.95	0.65
1:C:123:SER:OG	1:C:138:ASN:HB3	1.95	0.65
1:C:73:PHE:CE1	1:C:77:VAL:HG23	2.31	0.65
1:B:102:ARG:O	1:B:106:ILE:HD13	1.97	0.65
1:B:22:GLU:O	1:B:22:GLU:CG	2.45	0.65
1:C:107:ASN:O	1:C:110:PHE:HB3	1.97	0.64
1:B:53:ASP:OD1	1:B:60:THR:HG22	1.97	0.64
1:D:157:THR:HG22	1:D:166:ALA:HB2	1.80	0.64
1:D:24:TYR:CE2	1:D:33:ILE:HD11	2.33	0.64
1:C:112:MET:O	1:C:113:GLY:O	2.15	0.64
1:A:56:ILE:HD12	1:A:68:GLU:HB3	1.78	0.64
1:A:60:THR:OG1	1:A:63:VAL:O	2.16	0.64
1:A:14:ARG:HD3	1:A:18:TYR:CD1	2.32	0.63
1:D:19:LYS:NZ	1:D:43:PRO:HB2	2.13	0.63
1:A:18:TYR:O	1:A:25:TYR:HA	1.99	0.63
1:A:152:ALA:O	1:A:156:ILE:HG13	1.99	0.63
1:B:46:ASN:O	1:B:49:LYS:HB3	1.99	0.62
1:D:58:ARG:NH1	1:D:60:THR:HA	2.14	0.62
1:D:32:THR:O	1:D:33:ILE:CG1	2.44	0.62
1:C:144:TRP:CZ3	1:C:156:ILE:HD11	2.33	0.62
1:B:165:ASP:HA	1:B:168:LYS:NZ	2.13	0.62

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:4:PHE:HD2	1:C:73:PHE:CE2	2.18	0.62
1:A:69:ALA:O	1:A:72:LEU:HB2	1.99	0.62
1:C:3:ILE:O	1:C:6:MET:HB3	1.99	0.62
1:B:21:THR:CG2	1:B:21:THR:OG1	2.46	0.62
1:A:106:ILE:HG22	1:A:106:ILE:O	1.99	0.62
1:C:14:ARG:HH21	1:C:19:LYS:HB2	1.65	0.62
1:D:123:SER:O	1:D:127:LEU:HG	2.00	0.62
1:C:79:ALA:HA	1:C:82:ARG:NH1	2.15	0.62
1:C:106:ILE:HG22	1:C:106:ILE:O	2.00	0.61
1:D:17:ILE:CG2	1:D:17:ILE:C	2.68	0.61
1:C:25:TYR:OH	1:C:45:LEU:HG	1.99	0.61
1:A:78:ASP:O	1:A:81:VAL:HB	2.00	0.61
1:D:151:ARG:NH1	1:D:151:ARG:HG3	2.16	0.61
1:C:3:ILE:HG23	1:C:4:PHE:N	2.16	0.61
1:C:100:VAL:C	1:C:100:VAL:CG1	2.69	0.61
1:C:126:MET:SD	1:C:138:ASN:ND2	2.74	0.61
1:A:150:ASN:N	2:A:807:SO4:O4	2.33	0.60
1:B:161:THR:OG1	1:B:162:GLY:N	2.32	0.60
1:A:14:ARG:NH1	1:A:18:TYR:CE1	2.68	0.60
1:B:19:LYS:CG	1:B:19:LYS:H	2.14	0.60
1:D:13:LEU:HD11	1:D:15:LEU:HG	1.83	0.60
1:D:122:ASN:HB2	2:D:802:SO4:S	2.41	0.59
1:C:58:ARG:HH12	1:C:60:THR:HB	1.67	0.59
1:A:7:LEU:HD21	1:A:107:ASN:HA	1.83	0.59
1:B:148:THR:N	1:B:149:PRO:CD	2.65	0.59
1:D:13:LEU:HA	1:D:28:GLY:O	2.01	0.59
1:A:106:ILE:CG2	1:A:106:ILE:C	2.70	0.59
1:A:87:ASN:OD1	1:A:89:LYS:N	2.36	0.59
1:D:129:GLN:HB2	1:D:131:ARG:HD3	1.84	0.59
1:A:56:ILE:CD1	1:A:68:GLU:HB3	2.32	0.59
1:B:93:VAL:HG12	1:B:94:TYR:N	2.18	0.59
1:D:58:ARG:HH11	1:D:58:ARG:HG3	1.67	0.59
1:D:33:ILE:HG23	1:D:51:GLU:CG	2.28	0.59
1:B:87:ASN:OD1	1:B:89:LYS:N	2.36	0.59
1:B:150:ASN:ND2	1:B:150:ASN:H	1.99	0.59
1:C:145:TYR:O	1:C:149:PRO:HD3	2.03	0.59
1:A:52:LEU:HD23	1:A:60:THR:CG2	2.32	0.59
1:D:58:ARG:NH1	1:D:58:ARG:HG3	2.17	0.58
1:A:52:LEU:HD23	1:A:60:THR:HG21	1.84	0.58
1:D:17:ILE:HG22	1:D:18:TYR:N	2.17	0.58
1:B:138:ASN:HD22	1:B:141:LYS:HE2	1.67	0.58
1:D:3:ILE:HA	1:D:6:MET:HB3	1.84	0.58

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:29:ILE:HG21	1:A:73:PHE:CE1	2.38	0.58
1:D:45:LEU:O	1:D:49:LYS:HG3	2.04	0.58
1:C:64:ILE:HA	1:C:68:GLU:OE1	2.03	0.58
1:D:102:ARG:O	1:D:106:ILE:HG13	2.03	0.58
1:B:26:THR:HG23	1:B:31:TYR:O	2.03	0.58
1:B:90:LEU:O	1:B:93:VAL:HB	2.03	0.58
1:C:114:GLU:HG2	1:C:115:THR:N	2.18	0.58
1:C:59:ASN:ND2	1:C:61:ASN:HD21	2.02	0.57
1:B:19:LYS:CG	1:B:19:LYS:N	2.67	0.57
1:C:14:ARG:HE	1:C:19:LYS:CB	2.10	0.57
1:B:129:GLN:O	1:B:130:LYS:HB2	2.04	0.57
1:D:27:ILE:CG2	1:D:28:GLY:N	2.67	0.57
1:A:69:ALA:HA	1:A:72:LEU:HD12	1.87	0.57
1:C:17:ILE:HG22	1:C:18:TYR:N	2.19	0.57
1:A:106:ILE:N	1:A:106:ILE:CD1	2.66	0.57
1:A:81:VAL:O	1:A:84:ILE:HB	2.04	0.57
1:B:81:VAL:O	1:B:84:ILE:HB	2.04	0.57
1:B:16:LYS:O	1:B:17:ILE:HG23	2.04	0.57
1:C:122:ASN:HB2	2:C:806:SO4:O2	2.05	0.57
1:B:161:THR:OG1	1:B:163:THR:N	2.37	0.57
1:C:131:ARG:HH21	1:C:134:GLU:CD	2.09	0.57
1:A:142:SER:OG	1:A:144:TRP:HB3	2.05	0.56
1:B:2:ASN:OD1	1:B:5:GLU:N	2.29	0.56
1:D:13:LEU:HD13	1:D:13:LEU:C	2.26	0.56
1:A:106:ILE:HG22	1:A:106:ILE:C	2.25	0.56
1:C:14:ARG:NH2	1:C:19:LYS:HB2	2.20	0.56
1:B:138:ASN:HD22	1:B:141:LYS:CE	2.18	0.56
1:B:151:ARG:HA	1:B:154:ARG:NH2	2.17	0.56
1:C:93:VAL:HG11	1:C:124:LEU:HD23	1.87	0.56
1:B:82:ARG:HB2	1:B:82:ARG:HH11	1.69	0.56
1:D:7:LEU:HD23	1:D:107:ASN:ND2	2.21	0.56
1:C:90:LEU:HD23	1:C:124:LEU:HD11	1.88	0.56
1:B:150:ASN:O	1:B:153:LYS:HB2	2.06	0.56
1:D:19:LYS:HZ2	1:D:43:PRO:HB2	1.71	0.56
1:C:112:MET:HG3	1:C:117:VAL:HG22	1.87	0.56
1:A:51:GLU:O	1:A:54:LYS:HG3	2.06	0.56
1:A:19:LYS:HB3	1:A:25:TYR:CE2	2.40	0.56
1:C:86:ARG:CB	1:C:86:ARG:CZ	2.83	0.56
1:B:53:ASP:OD1	1:B:59:ASN:HA	2.05	0.56
1:C:93:VAL:HG21	1:C:124:LEU:HD22	1.88	0.56
1:D:13:LEU:HD13	1:D:14:ARG:N	2.21	0.55
1:D:60:THR:O	1:D:61:ASN:HB2	2.06	0.55

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:72:LEU:O	1:C:75:GLN:N	2.39	0.55
1:A:97:LEU:O	1:A:102:ARG:NE	2.40	0.55
1:A:120:PHE:CD2	1:A:120:PHE:N	2.74	0.55
1:C:109:VAL:HA	1:C:117:VAL:HG21	1.88	0.55
1:C:117:VAL:CG1	1:C:117:VAL:O	2.53	0.55
1:B:19:LYS:HG2	1:B:19:LYS:N	2.20	0.55
1:B:98:ASP:OD2	1:B:100:VAL:HG12	2.06	0.55
1:C:163:THR:HB	1:C:165:ASP:OD1	2.06	0.55
1:C:14:ARG:CZ	1:C:19:LYS:HB2	2.37	0.55
1:B:151:ARG:HA	1:B:154:ARG:HE	1.72	0.55
1:D:24:TYR:CB	1:D:24:TYR:C	2.71	0.55
1:B:100:VAL:HG11	1:B:162:GLY:C	2.27	0.55
1:D:165:ASP:HA	1:D:168:LYS:CE	2.36	0.55
1:C:143:ARG:HD3	1:C:143:ARG:O	2.06	0.54
1:B:148:THR:N	1:B:149:PRO:HD3	2.22	0.54
1:C:71:LYS:CE	1:C:75:GLN:HE21	2.06	0.54
1:D:24:TYR:CD2	1:D:33:ILE:HD11	2.43	0.54
1:B:84:ILE:HG21	1:B:94:TYR:CD1	2.43	0.54
1:D:157:THR:O	1:D:161:THR:HG23	2.07	0.54
1:C:124:LEU:HA	1:C:127:LEU:HD12	1.90	0.54
1:D:17:ILE:HG12	1:D:27:ILE:HD11	1.90	0.54
1:A:106:ILE:H	1:A:106:ILE:CD1	2.17	0.54
1:D:43:PRO:N	1:D:43:PRO:CG	2.69	0.53
1:C:91:LYS:HB3	1:C:92:PRO:HD3	1.89	0.53
1:A:27:ILE:HG22	1:A:28:GLY:H	1.73	0.53
1:B:131:ARG:HD3	1:B:134:GLU:HG3	1.89	0.53
1:C:144:TRP:HZ3	1:C:156:ILE:HD11	1.73	0.53
1:D:10:ASP:OD1	1:D:154:ARG:NH2	2.41	0.53
1:C:112:MET:O	1:C:117:VAL:HG23	2.08	0.53
1:B:165:ASP:HA	1:B:168:LYS:HZ1	1.72	0.53
1:A:15:LEU:O	1:A:63:VAL:HA	2.09	0.53
1:A:17:ILE:HG22	1:A:18:TYR:N	2.22	0.53
1:C:13:LEU:HD11	1:C:15:LEU:HD11	1.91	0.52
1:B:133:ASP:O	1:B:136:ALA:HB3	2.10	0.52
1:D:27:ILE:HG22	1:D:28:GLY:N	2.23	0.52
1:C:100:VAL:C	1:C:100:VAL:HG12	2.30	0.52
1:A:13:LEU:HD12	1:A:14:ARG:N	2.25	0.52
1:B:107:ASN:ND2	1:B:155:VAL:CG2	2.73	0.52
1:A:52:LEU:HD21	1:A:64:ILE:HG23	1.91	0.52
1:C:2:ASN:OD1	1:C:4:PHE:HB2	2.09	0.52
1:C:143:ARG:N	2:C:804:SO4:O1	2.42	0.52
1:B:147:GLN:O	1:B:147:GLN:HG3	2.10	0.52

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:74:ASN:ND2	1:B:78:ASP:OD1	2.42	0.52
1:D:98:ASP:OD2	1:D:100:VAL:HB	2.10	0.52
1:A:4:PHE:O	1:A:5:GLU:C	2.47	0.52
1:D:43:PRO:CB	1:D:43:PRO:CG	2.88	0.51
1:C:64:ILE:HG22	1:C:68:GLU:OE1	2.09	0.51
1:A:93:VAL:HG21	1:A:124:LEU:HB3	1.91	0.51
1:D:109:VAL:HA	1:D:117:VAL:CG2	2.40	0.51
1:D:17:ILE:HG23	1:D:18:TYR:H	1.75	0.51
1:A:120:PHE:HD2	1:A:120:PHE:N	2.07	0.51
1:A:2:ASN:OD1	1:A:4:PHE:HB2	2.11	0.51
1:A:126:MET:HE3	1:A:138:ASN:HD22	1.75	0.51
1:C:71:LYS:HG3	1:C:75:GLN:NE2	2.25	0.51
1:D:149:PRO:O	1:D:150:ASN:C	2.49	0.51
1:B:27:ILE:CD1	1:B:52:LEU:HD23	2.41	0.51
1:A:151:ARG:HB2	1:A:154:ARG:HH21	1.76	0.50
1:D:17:ILE:HG23	1:D:18:TYR:N	2.25	0.50
1:D:87:ASN:CB	1:D:114:GLU:OE2	2.48	0.50
1:A:143:ARG:HA	1:A:146:ASN:OD1	2.10	0.50
1:D:67:ASP:O	1:D:70:GLU:HB2	2.10	0.50
1:D:49:LYS:O	1:D:53:ASP:OD2	2.30	0.50
1:D:52:LEU:O	1:D:56:ILE:HG12	2.12	0.50
1:B:100:VAL:HG11	1:B:162:GLY:O	2.11	0.50
1:A:17:ILE:CG2	1:A:18:TYR:N	2.75	0.50
1:D:17:ILE:HA	1:D:27:ILE:HG13	1.93	0.50
1:B:107:ASN:HD22	1:B:155:VAL:HG21	1.77	0.50
1:B:49:LYS:HG3	1:B:53:ASP:OD2	2.10	0.50
1:A:53:ASP:HA	1:A:58:ARG:O	2.12	0.50
1:A:16:LYS:HG3	1:A:63:VAL:HG22	1.94	0.50
1:A:60:THR:HG22	1:A:60:THR:N	2.27	0.50
1:A:145:TYR:O	1:A:149:PRO:HD3	2.11	0.50
1:A:76:ASP:O	1:A:79:ALA:N	2.45	0.50
1:C:126:MET:SD	1:C:134:GLU:O	2.70	0.50
1:D:78:ASP:O	1:D:81:VAL:HB	2.12	0.50
1:D:26:THR:C	1:D:27:ILE:CD1	2.50	0.49
1:C:31:TYR:CD1	1:C:72:LEU:HD13	2.47	0.49
1:A:6:MET:SD	1:A:107:ASN:HB2	2.51	0.49
1:C:52:LEU:O	1:C:56:ILE:HG22	2.12	0.49
1:C:66:LYS:O	1:C:70:GLU:HG3	2.13	0.49
1:D:25:TYR:OH	1:D:45:LEU:HD22	2.13	0.49
1:C:65:THR:O	1:C:68:GLU:N	2.46	0.49
1:D:73:PHE:CD1	1:D:73:PHE:O	2.66	0.49
1:C:94:TYR:CE1	1:C:102:ARG:HB2	2.47	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:43:PRO:CG	1:D:43:PRO:CA	2.91	0.49
1:C:86:ARG:CB	1:C:86:ARG:NH1	2.73	0.49
1:C:16:LYS:HB3	1:C:16:LYS:NZ	2.28	0.49
1:B:53:ASP:O	1:B:56:ILE:HG22	2.12	0.49
1:C:1:MET:HG2	1:C:2:ASN:N	2.28	0.49
1:A:122:ASN:HB3	1:A:138:ASN:ND2	2.28	0.49
1:C:93:VAL:O	1:C:96:SER:HB3	2.12	0.49
1:D:105:LEU:O	1:D:109:VAL:HG23	2.12	0.49
1:D:91:LYS:N	1:D:92:PRO:CD	2.76	0.48
1:B:113:GLY:O	1:B:114:GLU:C	2.52	0.48
1:B:126:MET:HB3	1:B:135:ALA:HB2	1.95	0.48
1:A:126:MET:CE	1:A:138:ASN:ND2	2.76	0.48
1:B:72:LEU:O	1:B:75:GLN:HB2	2.12	0.48
1:D:17:ILE:CG2	1:D:17:ILE:CD1	2.91	0.48
1:D:52:LEU:HD12	1:D:56:ILE:CG1	2.43	0.48
1:C:27:ILE:HG13	1:C:64:ILE:CD1	2.44	0.48
1:A:68:GLU:O	1:A:72:LEU:HG	2.14	0.48
1:B:151:ARG:N	1:B:154:ARG:HH21	2.11	0.48
1:C:27:ILE:HG22	1:C:31:TYR:O	2.14	0.48
1:A:107:ASN:HD22	1:A:107:ASN:C	2.16	0.48
1:B:93:VAL:O	1:B:96:SER:HB3	2.14	0.48
1:C:152:ALA:O	1:C:156:ILE:HD13	2.14	0.48
1:C:59:ASN:ND2	1:C:61:ASN:ND2	2.62	0.48
1:C:135:ALA:O	1:C:136:ALA:C	2.50	0.48
1:A:11:GLU:HB3	1:A:30:GLY:H	1.79	0.47
1:D:125:ARG:O	1:D:128:GLN:HB3	2.14	0.47
1:D:58:ARG:HH12	1:D:60:THR:HA	1.79	0.47
1:C:7:LEU:HG	1:C:107:ASN:ND2	2.29	0.47
1:C:122:ASN:O	1:C:125:ARG:HG2	2.14	0.47
1:B:94:TYR:CE2	1:B:102:ARG:HD3	2.49	0.47
1:B:2:ASN:OD1	1:B:2:ASN:C	2.52	0.47
1:D:17:ILE:HG22	1:D:17:ILE:C	2.34	0.47
1:C:55:ALA:O	1:C:56:ILE:C	2.52	0.47
1:D:126:MET:HG2	1:D:131:ARG:HH21	1.79	0.47
1:D:105:LEU:HD12	1:D:105:LEU:HA	1.73	0.47
1:B:101:ARG:O	1:B:104:ALA:HB3	2.15	0.47
1:D:3:ILE:HD12	1:D:106:ILE:HB	1.97	0.47
1:C:52:LEU:HD12	1:C:56:ILE:CG2	2.39	0.47
1:D:82:ARG:O	1:D:86:ARG:HG3	2.15	0.47
1:A:26:THR:HB	1:A:31:TYR:O	2.15	0.47
1:D:126:MET:HA	1:D:131:ARG:HE	1.78	0.47
1:B:122:ASN:O	1:B:125:ARG:N	2.48	0.47

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:126:MET:SD	1:A:134:GLU:HB3	2.54	0.47
1:D:20:ASP:C	1:D:22:GLU:N	2.67	0.47
1:A:64:ILE:HD12	1:A:69:ALA:HA	1.97	0.47
1:C:84:ILE:HD11	1:C:109:VAL:HG21	1.97	0.47
1:C:127:LEU:O	1:C:128:GLN:C	2.50	0.47
1:C:82:ARG:O	1:C:85:LEU:N	2.47	0.47
1:B:104:ALA:O	1:B:108:MET:HG3	2.15	0.47
1:A:1:MET:CE	1:A:167:TYR:HB3	2.45	0.47
1:A:22:GLU:HG3	1:A:24:TYR:CE2	2.48	0.47
1:D:143:ARG:HG2	2:D:801:SO4:O1	2.15	0.47
1:C:73:PHE:CE1	1:C:77:VAL:CG2	2.98	0.46
1:C:53:ASP:OD1	1:C:60:THR:HG22	2.16	0.46
1:C:10:ASP:CG	1:C:151:ARG:HG3	2.34	0.46
1:C:96:SER:OG	1:C:96:SER:O	2.33	0.46
1:D:129:GLN:HB2	1:D:131:ARG:CD	2.46	0.46
1:A:101:ARG:CD	1:A:132:TRP:CH2	2.99	0.46
1:D:33:ILE:CG2	1:D:51:GLU:HG2	2.38	0.46
1:C:143:ARG:CD	1:C:143:ARG:O	2.64	0.46
1:D:18:TYR:CE2	1:D:20:ASP:HB3	2.50	0.46
1:D:18:TYR:O	1:D:25:TYR:HA	2.15	0.46
1:A:104:ALA:O	1:A:108:MET:HG3	2.16	0.46
1:C:126:MET:O	1:C:131:ARG:N	2.44	0.46
1:A:167:TYR:O	1:A:168:LYS:HG2	2.15	0.46
1:D:44:SER:O	1:D:47:ALA:HB3	2.15	0.46
1:C:5:GLU:O	1:C:9:ILE:CD1	2.60	0.46
1:D:158:THR:HG22	1:D:158:THR:O	2.16	0.46
1:A:14:ARG:O	1:A:28:GLY:N	2.46	0.46
1:A:66:LYS:HE2	1:A:67:ASP:OD1	2.15	0.46
1:A:109:VAL:O	1:A:113:GLY:N	2.42	0.46
1:B:133:ASP:O	1:B:136:ALA:N	2.49	0.46
1:C:13:LEU:CD1	1:C:15:LEU:CD1	2.94	0.46
1:B:93:VAL:O	1:B:94:TYR:C	2.54	0.46
1:C:109:VAL:HG13	1:C:114:GLU:HA	1.97	0.46
1:B:72:LEU:HA	1:B:75:GLN:NE2	2.30	0.46
1:C:3:ILE:CG2	1:C:4:PHE:N	2.78	0.45
1:C:74:ASN:O	1:C:77:VAL:HB	2.16	0.45
1:C:82:ARG:O	1:C:83:GLY:C	2.53	0.45
1:A:101:ARG:NE	1:A:132:TRP:CZ2	2.84	0.45
1:A:109:VAL:HG13	1:A:114:GLU:N	2.31	0.45
1:B:142:SER:CB	1:B:144:TRP:HB3	2.47	0.45
1:B:10:ASP:HB3	1:B:151:ARG:CD	2.42	0.45
1:D:7:LEU:CD2	1:D:107:ASN:CG	2.85	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:142:SER:CB	1:B:144:TRP:H	2.29	0.45
1:A:117:VAL:C	1:A:119:GLY:N	2.70	0.45
1:A:85:LEU:N	1:A:85:LEU:HD12	2.30	0.45
1:D:52:LEU:HD12	1:D:56:ILE:HG12	1.97	0.45
1:C:99:ALA:HA	1:C:102:ARG:HG2	1.98	0.45
1:A:29:ILE:CG2	1:A:73:PHE:CE1	2.99	0.45
1:C:48:ALA:O	1:C:51:GLU:N	2.50	0.45
1:B:93:VAL:O	1:B:96:SER:N	2.48	0.45
1:C:163:THR:CB	1:C:165:ASP:OD1	2.64	0.45
1:B:113:GLY:O	1:B:116:GLY:N	2.49	0.45
1:C:21:THR:OG1	1:C:22:GLU:N	2.48	0.45
1:A:14:ARG:NH1	1:A:18:TYR:CZ	2.85	0.45
1:D:7:LEU:HD21	1:D:107:ASN:CG	2.36	0.45
1:D:150:ASN:O	1:D:154:ARG:HG3	2.16	0.45
1:D:2:ASN:O	1:D:6:MET:N	2.40	0.45
1:A:31:TYR:CD2	1:A:31:TYR:C	2.90	0.45
1:A:101:ARG:CZ	1:A:132:TRP:CZ2	2.99	0.45
1:A:150:ASN:O	1:A:154:ARG:HG3	2.17	0.45
1:C:3:ILE:HG23	1:C:4:PHE:H	1.82	0.45
1:C:165:ASP:HA	1:C:168:LYS:HG2	1.97	0.45
1:D:66:LYS:O	1:D:70:GLU:HG3	2.16	0.45
1:D:58:ARG:CG	1:D:58:ARG:HH11	2.28	0.44
1:C:14:ARG:HH21	1:C:19:LYS:CB	2.27	0.44
1:D:18:TYR:HE2	1:D:20:ASP:HB3	1.82	0.44
1:D:49:LYS:HA	1:D:52:LEU:HB3	1.98	0.44
1:A:124:LEU:O	1:A:125:ARG:C	2.54	0.44
1:C:80:ALA:O	1:C:81:VAL:C	2.53	0.44
1:D:27:ILE:HG21	1:D:64:ILE:HD12	1.98	0.44
1:A:120:PHE:HB3	1:A:123:SER:HB2	1.99	0.44
1:C:140:ALA:HA	1:C:145:TYR:CD1	2.52	0.44
1:B:64:ILE:CD1	1:B:69:ALA:HB2	2.48	0.44
1:B:138:ASN:HA	1:B:141:LYS:HE2	1.99	0.44
1:C:81:VAL:O	1:C:82:ARG:C	2.56	0.44
1:D:86:ARG:HB2	1:D:86:ARG:CZ	2.48	0.44
1:C:26:THR:O	1:C:27:ILE:HB	2.18	0.44
1:A:126:MET:HE3	1:A:138:ASN:ND2	2.33	0.44
1:D:150:ASN:O	1:D:151:ARG:C	2.55	0.44
1:C:126:MET:HB3	1:C:131:ARG:HB2	1.99	0.44
1:B:72:LEU:HD23	1:B:75:GLN:NE2	2.33	0.44
1:D:58:ARG:CZ	1:D:60:THR:HA	2.48	0.44
1:B:26:THR:HG22	1:B:27:ILE:N	2.32	0.44
1:A:100:VAL:HG22	1:A:164:TRP:CE2	2.52	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:98:ASP:O	1:B:102:ARG:N	2.44	0.43
1:B:17:ILE:HD12	1:B:17:ILE:O	2.18	0.43
1:D:129:GLN:O	1:D:130:LYS:C	2.56	0.43
1:D:45:LEU:CD2	1:D:45:LEU:H	2.03	0.43
1:C:71:LYS:HE3	1:C:71:LYS:HA	1.99	0.43
1:C:72:LEU:HD23	1:C:72:LEU:HA	1.67	0.43
1:B:16:LYS:HG3	1:B:17:ILE:N	2.33	0.43
1:B:27:ILE:HD12	1:B:52:LEU:HD23	2.00	0.43
1:A:14:ARG:NH1	1:A:18:TYR:CD1	2.86	0.43
1:D:98:ASP:OD2	1:D:101:ARG:HG3	2.18	0.43
1:A:98:ASP:C	1:A:102:ARG:HG2	2.35	0.43
1:B:53:ASP:CA	1:B:56:ILE:HG22	2.49	0.43
1:D:12:GLY:O	1:D:29:ILE:HG23	2.18	0.43
1:C:130:LYS:HA	1:C:132:TRP:CH2	2.53	0.43
1:A:7:LEU:O	1:A:11:GLU:N	2.52	0.43
1:B:53:ASP:C	1:B:56:ILE:HG22	2.39	0.43
1:D:154:ARG:O	1:D:158:THR:OG1	2.20	0.43
1:A:51:GLU:HA	1:A:54:LYS:HG2	2.00	0.43
1:A:117:VAL:C	1:A:119:GLY:H	2.22	0.43
1:B:13:LEU:HD12	1:B:14:ARG:H	1.83	0.43
1:D:3:ILE:CD1	1:D:106:ILE:CG2	2.97	0.42
1:A:2:ASN:OD1	1:A:2:ASN:C	2.56	0.42
1:C:65:THR:OG1	1:C:67:ASP:HB2	2.18	0.42
1:D:91:LYS:CB	1:D:92:PRO:HD3	2.42	0.42
1:A:76:ASP:O	1:A:79:ALA:CB	2.64	0.42
1:A:81:VAL:HG12	1:A:82:ARG:N	2.34	0.42
1:B:22:GLU:HG3	1:B:23:GLY:N	2.32	0.42
1:D:150:ASN:HA	1:D:153:LYS:HD2	2.01	0.42
1:D:7:LEU:HD23	1:D:7:LEU:HA	1.86	0.42
1:A:66:LYS:HD2	1:A:67:ASP:N	2.34	0.42
1:A:19:LYS:O	1:A:19:LYS:HD3	2.19	0.42
1:D:132:TRP:HA	1:D:132:TRP:CE3	2.54	0.42
1:A:131:ARG:O	1:A:132:TRP:C	2.58	0.42
1:D:151:ARG:NH1	1:D:151:ARG:CG	2.83	0.42
1:C:11:GLU:HB2	1:C:29:ILE:CG2	2.49	0.42
1:D:123:SER:N	2:D:802:SO4:O3	2.49	0.42
1:C:91:LYS:O	1:C:94:TYR:HB3	2.19	0.42
1:B:71:LYS:O	1:B:75:GLN:HG3	2.19	0.42
1:D:52:LEU:CD1	1:D:56:ILE:CG1	2.98	0.42
1:D:64:ILE:HD13	1:D:64:ILE:N	2.27	0.42
1:A:11:GLU:CD	1:A:151:ARG:HH12	2.22	0.42
1:D:17:ILE:HG21	1:D:17:ILE:CD1	2.49	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:16:LYS:O	1:C:17:ILE:C	2.58	0.42
1:B:16:LYS:O	1:B:27:ILE:CG2	2.68	0.42
1:D:84:ILE:C	1:D:86:ARG:H	2.22	0.42
1:A:71:LYS:O	1:A:75:GLN:HG3	2.20	0.42
1:D:52:LEU:CG	1:D:60:THR:HG21	2.36	0.42
1:A:98:ASP:O	1:A:102:ARG:CG	2.54	0.42
1:B:53:ASP:OD1	1:B:60:THR:N	2.53	0.42
1:B:152:ALA:O	1:B:156:ILE:HG13	2.19	0.42
1:D:17:ILE:HA	1:D:27:ILE:CG1	2.50	0.41
1:C:11:GLU:O	1:C:12:GLY:C	2.58	0.41
1:A:27:ILE:N	1:A:27:ILE:HD12	2.34	0.41
1:A:82:ARG:C	1:A:84:ILE:H	2.24	0.41
1:A:113:GLY:O	1:A:117:VAL:HG23	2.20	0.41
1:A:102:ARG:O	1:A:105:LEU:HB3	2.20	0.41
1:B:26:THR:HA	1:B:31:TYR:O	2.20	0.41
1:C:13:LEU:HD11	1:C:15:LEU:HD12	2.01	0.41
1:D:131:ARG:O	1:D:132:TRP:C	2.58	0.41
1:D:84:ILE:O	1:D:86:ARG:N	2.53	0.41
1:A:91:LYS:O	1:A:92:PRO:C	2.53	0.41
1:C:1:MET:HG3	1:C:5:GLU:CB	2.38	0.41
1:A:108:MET:HE1	1:A:144:TRP:CZ3	2.56	0.41
1:B:147:GLN:O	1:B:147:GLN:CG	2.68	0.41
1:A:167:TYR:C	1:A:168:LYS:HG2	2.40	0.41
1:B:107:ASN:CG	1:B:151:ARG:HH21	2.24	0.41
1:B:52:LEU:HD12	1:B:60:THR:HG21	2.02	0.41
1:D:130:LYS:HD2	1:D:132:TRP:CH2	2.55	0.41
1:D:97:LEU:HD22	1:D:101:ARG:CB	2.51	0.41
1:B:3:ILE:HG23	1:B:4:PHE:CD1	2.56	0.41
1:D:165:ASP:CA	1:D:168:LYS:HE2	2.46	0.41
1:A:52:LEU:HD23	1:A:60:THR:HG23	2.01	0.41
1:B:82:ARG:CB	1:B:82:ARG:HH11	2.34	0.41
1:A:19:LYS:CB	1:A:25:TYR:CE2	3.03	0.41
1:D:82:ARG:O	1:D:83:GLY:C	2.58	0.41
1:A:60:THR:CG2	1:A:60:THR:H	2.31	0.41
1:B:27:ILE:HB	1:B:28:GLY:H	1.68	0.41
1:B:16:LYS:HG3	1:B:17:ILE:H	1.86	0.41
1:C:137:VAL:HG12	1:C:141:LYS:NZ	2.35	0.41
1:B:45:LEU:HD12	1:B:48:ALA:HB3	2.02	0.41
1:C:52:LEU:C	1:C:52:LEU:HD12	2.41	0.41
1:B:98:ASP:HA	1:B:102:ARG:HH21	1.86	0.41
1:A:124:LEU:C	1:A:126:MET:N	2.72	0.41
1:B:5:GLU:O	1:B:8:ARG:HB3	2.21	0.41

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:86:ARG:HB2	1:D:86:ARG:NH1	2.36	0.41
1:C:8:ARG:HH21	1:C:13:LEU:HD23	1.86	0.40
1:B:7:LEU:O	1:B:8:ARG:C	2.57	0.40
1:C:53:ASP:OD1	1:C:60:THR:N	2.54	0.40
1:A:120:PHE:O	1:A:124:LEU:HG	2.21	0.40
1:D:164:TRP:O	1:D:168:LYS:HG3	2.21	0.40
1:C:100:VAL:HG12	1:C:100:VAL:O	2.21	0.40
1:C:27:ILE:CG1	1:C:64:ILE:HD11	2.51	0.40
1:B:133:ASP:O	1:B:137:VAL:HG23	2.22	0.40
1:D:5:GLU:O	1:D:9:ILE:HG12	2.21	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	153/170 (90%)	124 (81%)	22 (14%)	7 (5%)	4	22
1	B	153/170 (90%)	119 (78%)	27 (18%)	7 (5%)	4	22
1	C	153/170 (90%)	112 (73%)	31 (20%)	10 (6%)	2	11
1	D	156/170 (92%)	119 (76%)	30 (19%)	7 (4%)	4	22
All	All	615/680 (90%)	474 (77%)	110 (18%)	31 (5%)	3	19

All (31) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	17	ILE
1	C	113	GLY
1	D	43	PRO
1	D	118	ALA
1	A	30	GLY
1	A	81	VAL
1	A	132	TRP
1	B	93	VAL

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Mol	Chain	Res	Type
1	C	20	ASP
1	C	144	TRP
1	D	149	PRO
1	A	13	LEU
1	A	144	TRP
1	B	17	ILE
1	B	150	ASN
1	B	165	ASP
1	C	21	THR
1	C	55	ALA
1	C	93	VAL
1	D	22	GLU
1	D	150	ASN
1	B	97	LEU
1	C	12	GLY
1	C	27	ILE
1	C	81	VAL
1	D	141	LYS
1	B	94	TYR
1	B	114	GLU
1	D	85	LEU
1	A	165	ASP
1	A	77	VAL

### 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	130/141 (92%)	107 (82%)	23 (18%)	3	14
1	B	130/141 (92%)	112 (86%)	18 (14%)	5	24
1	C	130/141 (92%)	111 (85%)	19 (15%)	5	21
1	D	133/141 (94%)	113 (85%)	20 (15%)	4	20
All	All	523/564 (93%)	443 (85%)	80 (15%)	4	19

All (80) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	TYR
1	A	19	LYS
1	A	22	GLU
1	A	24	TYR
1	A	25	TYR
1	A	27	ILE
1	A	61	ASN
1	A	66	LYS
1	A	67	ASP
1	A	78	ASP
1	A	102	ARG
1	A	107	ASN
1	A	114	GLU
1	A	115	THR
1	A	120	PHE
1	A	122	ASN
1	A	125	ARG
1	A	128	GLN
1	A	129	GLN
1	A	137	VAL
1	A	143	ARG
1	A	146	ASN
1	A	163	THR
1	B	20	ASP
1	B	22	GLU
1	B	27	ILE
1	B	52	LEU
1	B	65	THR
1	B	74	ASN
1	B	82	ARG
1	B	96	SER
1	B	100	VAL
1	B	121	THR
1	B	123	SER
1	B	130	LYS
1	B	131	ARG
1	B	138	ASN
1	B	143	ARG
1	B	150	ASN
1	B	165	ASP
1	B	168	LYS
1	C	8	ARG
1	C	9	ILE

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Mol	Chain	Res	Type
1	C	10	ASP
1	C	16	LYS
1	C	20	ASP
1	C	27	ILE
1	C	45	LEU
1	C	52	LEU
1	C	56	ILE
1	C	60	THR
1	C	64	ILE
1	C	65	THR
1	C	71	LYS
1	C	78	ASP
1	C	86	ARG
1	C	96	SER
1	C	124	LEU
1	C	143	ARG
1	C	156	ILE
1	D	17	ILE
1	D	32	THR
1	D	43	PRO
1	D	44	SER
1	D	46	ASN
1	D	52	LEU
1	D	56	ILE
1	D	60	THR
1	D	64	ILE
1	D	67	ASP
1	D	72	LEU
1	D	73	PHE
1	D	74	ASN
1	D	78	ASP
1	D	114	GLU
1	D	131	ARG
1	D	141	LYS
1	D	151	ARG
1	D	157	THR
1	D	161	THR

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	107	ASN

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Mol	Chain	Res	Type
1	A	129	GLN
1	A	147	GLN
1	B	74	ASN
1	B	75	GLN
1	B	107	ASN
1	B	138	ASN
1	B	146	ASN
1	B	150	ASN
1	C	59	ASN
1	C	75	GLN
1	D	74	ASN
1	D	147	GLN

### 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 ⓘ

8 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z  > 2$	Counts	RMSZ	$\# Z  > 2$
2	SO4	A	807	-	4,4,4	0.54	0	6,6,6	0.90	0
2	SO4	A	808	-	4,4,4	0.39	0	6,6,6	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	SO4	C	804	-	4,4,4	0.18	0	6,6,6	0.57	0
2	SO4	C	805	-	4,4,4	0.41	0	6,6,6	0.45	0
2	SO4	C	806	-	4,4,4	0.61	0	6,6,6	0.68	0
2	SO4	D	801	-	4,4,4	0.76	0	6,6,6	0.73	0
2	SO4	D	802	-	4,4,4	0.22	0	6,6,6	0.25	0
2	SO4	D	803	-	4,4,4	0.46	0	6,6,6	0.56	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
2	SO4	A	807	-	-	0/0/0/0	0/0/0/0
2	SO4	A	808	-	-	0/0/0/0	0/0/0/0
2	SO4	C	804	-	-	0/0/0/0	0/0/0/0
2	SO4	C	805	-	-	0/0/0/0	0/0/0/0
2	SO4	C	806	-	-	0/0/0/0	0/0/0/0
2	SO4	D	801	-	-	0/0/0/0	0/0/0/0
2	SO4	D	802	-	-	0/0/0/0	0/0/0/0
2	SO4	D	803	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	157/170 (92%)	-0.54	0 100 100	2, 7, 15, 20	0
1	B	157/170 (92%)	-0.55	0 100 100	1, 7, 19, 36	0
1	C	157/170 (92%)	-0.57	0 100 100	2, 7, 20, 37	0
1	D	160/170 (94%)	-0.54	0 100 100	1, 6, 24, 35	0
All	All	631/680 (92%)	-0.55	0 100 100	1, 7, 19, 37	0

There are no RSRZ outliers to report.

### 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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
2	SO4	D	803	5/5	0.14	0.47	37,37,39,41	0
2	SO4	C	806	5/5	0.14	0.03	29,33,34,36	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SO4	C	805	5/5	0.13	-0.75	21,25,26,33	0
2	SO4	D	801	5/5	0.12	-1.22	13,15,19,19	0
2	SO4	C	804	5/5	0.11	-1.43	14,14,22,26	0
2	SO4	A	808	5/5	0.10	-1.64	30,31,32,34	0
2	SO4	D	802	5/5	0.09	-1.65	37,38,41,42	0
2	SO4	A	807	5/5	0.09	-1.82	20,20,24,26	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.