



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:15 AM GMT

PDB ID : 2C44  
Title : CRYSTAL STRUCTURE OF E. COLI TRYPTOPHANASE  
Authors : Ku, S.-Y.; Yip, P.; Howell, P.L.  
Deposited on : 2005-10-15  
Resolution : 2.80 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](mailto:validation@mail.wwpdb.org)  
A user guide is available at  
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>  
with specific help available everywhere you see the ⓘ symbol.

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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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

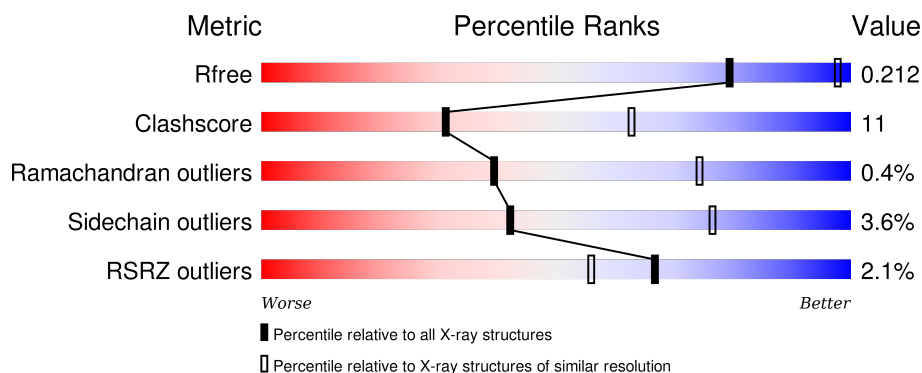
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 2.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
$R_{free}$	91344	2393 (2.80-2.80)
Clashscore	102246	2827 (2.80-2.80)
Ramachandran outliers	100387	2782 (2.80-2.80)
Sidechain outliers	100360	2784 (2.80-2.80)
RSRZ outliers	91569	2404 (2.80-2.80)

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. 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	471	<div> <div>7%</div> <div>78%19%...</div> </div>
1	B	471	<div> <div>79%18%...</div> </div>
1	C	471	<div> <div>78%19%...</div> </div>
1	D	471	<div> <div>7%</div> <div>78%18%...</div> </div>

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 14846 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 TRYPTOPHANASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	0	0	1
			3570	2278	599	671	22			
1	B	467	Total	C	N	O	S	0	0	0
			3584	2288	604	670	22			
1	C	468	Total	C	N	O	S	0	0	0
			3584	2288	605	669	22			
1	D	467	Total	C	N	O	S	0	0	1
			3544	2258	598	666	22			

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



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		

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

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	K	0	0
			1	1		
3	A	2	Total	K	0	0
			2	2		
3	D	1	Total	K	0	0
			1	1		
3	C	2	Total	K	0	0
			2	2		

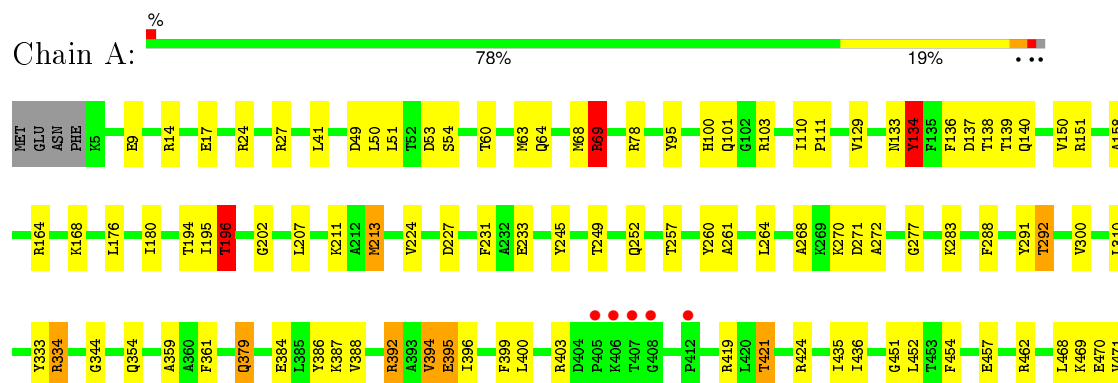
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	130	Total	O	0	0
			130	130		
4	B	143	Total	O	0	0
			143	143		
4	C	125	Total	O	0	0
			125	125		
4	D	120	Total	O	0	0
			120	120		

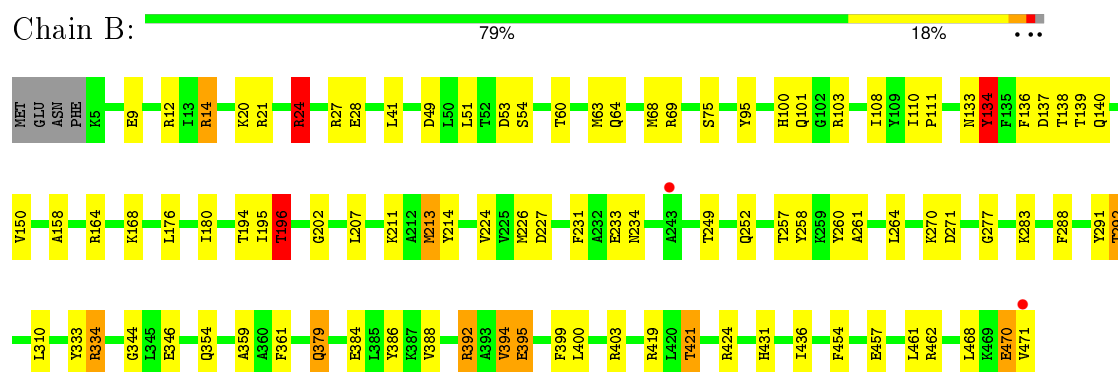
### 3 Residue-property plots

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

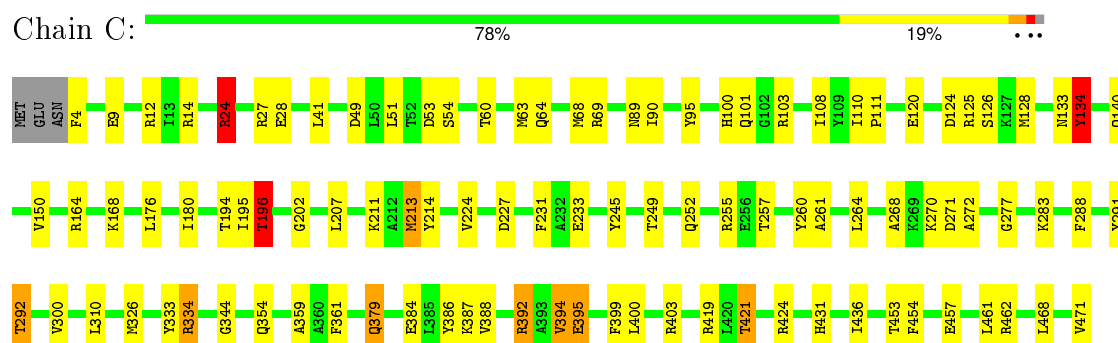
#### • Molecule 1: TRYPTOPHANASE



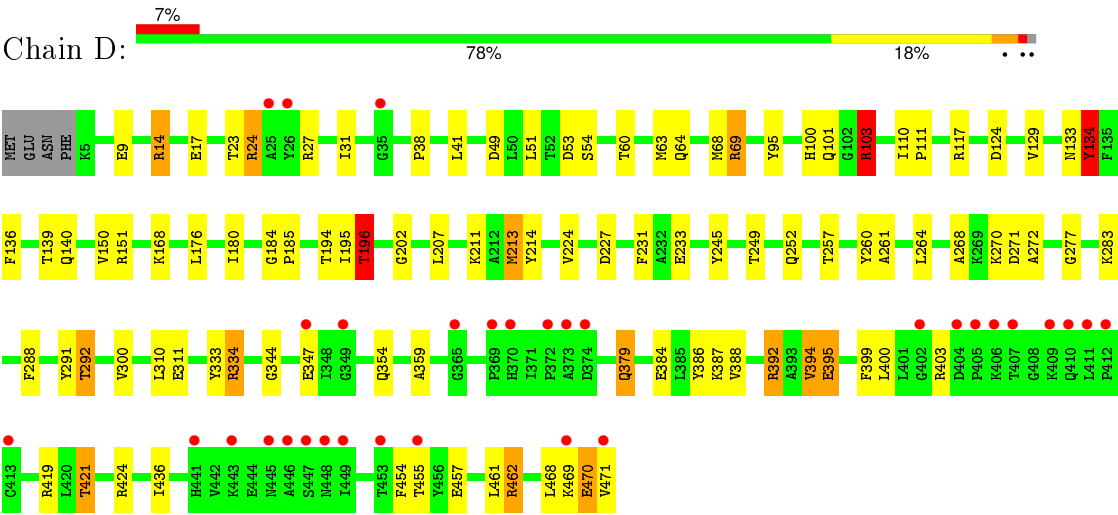
#### • Molecule 1: TRYPTOPHANASE



#### • Molecule 1: TRYPTOPHANASE



● Molecule 1: TRYPTOPHANASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	215.51Å 215.51Å 107.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	500.00 – 2.80 48.19 – 2.81	Depositor EDS
% Data completeness (in resolution range)	99.3 (500.00-2.80) 99.9 (48.19-2.81)	Depositor EDS
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	6.64 (at 2.81Å)	Xtriage
Refinement program	REFMAC 5.2.0005	Depositor
R, $R_{free}$	0.196 , 0.220 0.192 , 0.212	Depositor DCC
$R_{free}$ test set	3141 reflections (5.31%)	DCC
Wilson B-factor (Å <sup>2</sup> )	17.7	Xtriage
Anisotropy	0.275	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 38.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.45$ , $\langle L^2 \rangle = 0.28$	Xtriage
Outliers	0 of 62256 reflections	Xtriage
$F_o, F_c$ correlation	0.91	EDS
Total number of atoms	14846	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.25% of the height of the origin peak. No significant pseudotranslation is detected.*

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

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

## 5 Model quality ⓘ

### 5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: K, 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	0.36	0/3648	0.89	10/4952 (0.2%)
1	B	0.35	0/3662	0.68	8/4967 (0.2%)
1	C	0.34	0/3662	0.68	9/4968 (0.2%)
1	D	0.37	0/3622	0.86	10/4920 (0.2%)
All	All	0.36	0/14594	0.78	37/19807 (0.2%)

There are no bond length outliers.

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	69	ARG	NE-CZ-NH2	-21.55	109.53	120.30
1	A	24	ARG	NE-CZ-NH2	-20.63	109.98	120.30
1	A	69	ARG	NE-CZ-NH1	20.51	130.56	120.30
1	A	24	ARG	NE-CZ-NH1	19.85	130.23	120.30
1	D	462	ARG	NE-CZ-NH1	-19.17	110.72	120.30
1	D	462	ARG	NE-CZ-NH2	18.90	129.75	120.30
1	D	103	ARG	NE-CZ-NH2	-17.80	111.40	120.30
1	D	103	ARG	NE-CZ-NH1	17.23	128.91	120.30
1	A	69	ARG	CD-NE-CZ	10.59	138.43	123.60
1	A	24	ARG	CD-NE-CZ	10.15	137.82	123.60
1	D	462	ARG	CD-NE-CZ	8.85	135.99	123.60
1	D	69	ARG	NE-CZ-NH1	-7.54	116.53	120.30
1	C	69	ARG	NE-CZ-NH1	-7.52	116.54	120.30
1	C	69	ARG	NE-CZ-NH2	7.25	123.92	120.30
1	B	69	ARG	NE-CZ-NH2	7.23	123.92	120.30
1	C	24	ARG	NE-CZ-NH1	-7.22	116.69	120.30
1	D	24	ARG	NE-CZ-NH1	-7.09	116.76	120.30
1	D	69	ARG	NE-CZ-NH2	7.04	123.82	120.30
1	D	103	ARG	CD-NE-CZ	7.02	133.43	123.60
1	B	24	ARG	NE-CZ-NH1	-6.97	116.81	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	69	ARG	NE-CZ-NH1	-6.95	116.83	120.30
1	D	24	ARG	NE-CZ-NH2	6.85	123.73	120.30
1	C	24	ARG	NE-CZ-NH2	6.68	123.64	120.30
1	B	462	ARG	NE-CZ-NH2	-6.43	117.09	120.30
1	C	103	ARG	NE-CZ-NH1	-6.35	117.12	120.30
1	A	462	ARG	NE-CZ-NH2	-6.29	117.15	120.30
1	C	462	ARG	NE-CZ-NH2	-6.21	117.20	120.30
1	B	24	ARG	NE-CZ-NH2	6.06	123.33	120.30
1	B	103	ARG	NE-CZ-NH2	5.68	123.14	120.30
1	A	103	ARG	NE-CZ-NH1	-5.67	117.47	120.30
1	B	103	ARG	NE-CZ-NH1	-5.59	117.50	120.30
1	C	103	ARG	NE-CZ-NH2	5.50	123.05	120.30
1	C	462	ARG	NE-CZ-NH1	5.49	123.04	120.30
1	A	462	ARG	NE-CZ-NH1	5.40	123.00	120.30
1	B	462	ARG	NE-CZ-NH1	5.09	122.85	120.30
1	A	103	ARG	NE-CZ-NH2	5.06	122.83	120.30
1	C	196	THR	N-CA-C	-5.03	97.43	111.00

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	A	3570	0	3432	78	0
1	B	3584	0	3463	73	0
1	C	3584	0	3459	79	0
1	D	3544	0	3372	87	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
2	C	10	0	0	0	0
2	D	10	0	0	1	0
3	A	2	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	0	0
3	D	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	A	130	0	0	3	0
4	B	143	0	0	2	0
4	C	125	0	0	4	0
4	D	120	0	0	10	0
All	All	14846	0	13726	302	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 11.

All (302) 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:D:60:THR:H	1:D:63:MET:HE3	1.37	0.89
1:C:60:THR:H	1:C:63:MET:HE3	1.40	0.86
1:B:60:THR:H	1:B:63:MET:HE3	1.40	0.86
1:D:51:LEU:HD12	1:D:394:VAL:HG13	1.58	0.85
1:A:60:THR:H	1:A:63:MET:HE3	1.39	0.85
1:A:51:LEU:HD12	1:A:394:VAL:HG13	1.59	0.84
1:B:51:LEU:HD12	1:B:394:VAL:HG13	1.59	0.84
1:A:51:LEU:HB2	1:A:392:ARG:HD3	1.60	0.83
1:A:452:LEU:HD23	1:A:471:VAL:N	1.93	0.83
1:C:51:LEU:HD12	1:C:394:VAL:HG13	1.59	0.82
1:D:51:LEU:HB2	1:D:392:ARG:HD3	1.62	0.81
1:B:51:LEU:HB2	1:B:392:ARG:HD3	1.61	0.80
1:C:51:LEU:HB2	1:C:392:ARG:HD3	1.62	0.79
1:B:470:GLU:O	1:B:471:VAL:HG12	1.85	0.76
1:A:379:GLN:HG2	1:A:395:GLU:HG2	1.68	0.76
1:D:129:VAL:HG11	1:D:151:ARG:HD3	1.68	0.75
1:D:379:GLN:HG2	1:D:395:GLU:HG2	1.69	0.75
1:B:471:VAL:HG23	4:B:2061:HOH:O	1.87	0.75
1:C:379:GLN:HG2	1:C:395:GLU:HG2	1.70	0.73
1:B:379:GLN:HG2	1:B:395:GLU:HG2	1.70	0.72
1:A:68:MET:HE3	1:D:64:GLN:HG2	1.72	0.72
1:A:359:ALA:HB2	1:A:421:THR:HA	1.73	0.70
1:D:60:THR:N	1:D:63:MET:HE3	2.06	0.70
1:D:136:PHE:HB2	1:D:139:THR:OG1	1.92	0.70
1:D:359:ALA:HB2	1:D:421:THR:HA	1.74	0.70
1:D:424:ARG:HD3	4:D:2115:HOH:O	1.90	0.70
1:B:359:ALA:HB2	1:B:421:THR:HA	1.73	0.69
1:C:359:ALA:HB2	1:C:421:THR:HA	1.74	0.69
1:A:69:ARG:NH2	1:D:69:ARG:NH2	2.41	0.69

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:THR:N	1:A:63:MET:HE3	2.08	0.68
1:A:64:GLN:HG2	1:D:68:MET:HE3	1.75	0.68
1:B:60:THR:N	1:B:63:MET:HE3	2.09	0.67
1:D:103:ARG:HD2	2:D:500:SO4:O2	1.95	0.66
1:C:196:THR:HG21	4:C:2057:HOH:O	1.94	0.66
1:A:69:ARG:NH2	1:D:69:ARG:HH22	1.94	0.66
1:C:60:THR:N	1:C:63:MET:HE3	2.09	0.65
1:A:69:ARG:HH21	1:D:69:ARG:HH22	1.44	0.65
1:B:419:ARG:HD3	1:B:421:THR:HG23	1.78	0.65
1:D:419:ARG:HD3	1:D:421:THR:HG23	1.79	0.64
1:C:419:ARG:HD3	1:C:421:THR:HG23	1.80	0.64
1:D:38:PRO:HA	4:D:2013:HOH:O	1.98	0.63
1:A:419:ARG:HD3	1:A:421:THR:HG23	1.80	0.63
1:D:136:PHE:HE1	1:D:194:THR:HG23	1.64	0.62
1:C:195:ILE:O	1:C:196:THR:HB	2.00	0.62
1:C:164:ARG:HG2	1:C:164:ARG:HH21	1.65	0.61
1:B:195:ILE:O	1:B:196:THR:HB	2.00	0.60
1:B:64:GLN:HG2	1:C:68:MET:HE3	1.83	0.60
1:D:49:ASP:OD2	1:D:392:ARG:HD2	2.03	0.59
1:B:400:LEU:C	1:B:400:LEU:HD23	2.24	0.58
1:D:38:PRO:HG3	4:D:2013:HOH:O	2.01	0.58
1:D:195:ILE:HG21	1:D:231:PHE:HA	1.86	0.58
1:A:195:ILE:O	1:A:196:THR:HB	2.04	0.57
1:A:49:ASP:OD2	1:A:392:ARG:HD2	2.04	0.57
1:D:196:THR:HG21	4:D:2066:HOH:O	2.04	0.57
1:C:24:ARG:HD2	1:C:28:GLU:OE2	2.03	0.57
1:C:24:ARG:O	1:C:28:GLU:HG2	2.04	0.57
1:B:195:ILE:HG21	1:B:231:PHE:HA	1.85	0.57
1:B:24:ARG:HD2	1:B:28:GLU:OE1	2.04	0.57
1:C:49:ASP:OD2	1:C:392:ARG:HD2	2.05	0.57
1:B:471:VAL:OXT	1:B:471:VAL:HG13	2.04	0.57
1:C:195:ILE:HG21	1:C:231:PHE:HA	1.86	0.56
1:C:164:ARG:HG2	1:C:164:ARG:NH2	2.19	0.56
1:C:277:GLY:HA2	1:C:310:LEU:HD21	1.87	0.56
1:C:400:LEU:HD23	1:C:400:LEU:C	2.25	0.56
1:B:49:ASP:OD2	1:B:392:ARG:HD2	2.05	0.56
1:D:100:HIS:CD2	1:D:101:GLN:HG2	2.41	0.56
1:D:195:ILE:CG2	1:D:231:PHE:HA	2.36	0.55
1:C:53:ASP:OD2	1:C:424:ARG:NH2	2.39	0.55
1:D:53:ASP:OD2	1:D:424:ARG:NH2	2.40	0.55
1:A:195:ILE:HG21	1:A:231:PHE:HA	1.87	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:ILE:CG2	1:B:231:PHE:HA	2.37	0.55
1:A:110:ILE:HB	1:A:111:PRO:HD3	1.89	0.55
1:B:54:SER:HB2	1:B:270:LYS:HE3	1.89	0.55
1:A:400:LEU:C	1:A:400:LEU:HD23	2.26	0.55
1:D:277:GLY:HA2	1:D:310:LEU:HD21	1.89	0.55
1:B:100:HIS:CD2	1:B:101:GLN:HG2	2.42	0.55
1:D:195:ILE:O	1:D:196:THR:HB	2.05	0.54
1:D:110:ILE:HB	1:D:111:PRO:HD3	1.90	0.54
1:A:133:ASN:O	1:A:134:TYR:HB2	2.07	0.54
1:A:54:SER:HB2	1:A:270:LYS:HE3	1.90	0.54
1:D:400:LEU:HD23	1:D:400:LEU:C	2.27	0.54
1:C:195:ILE:CG2	1:C:231:PHE:HA	2.37	0.54
1:D:264:LEU:HD12	1:D:264:LEU:C	2.28	0.54
1:B:110:ILE:HB	1:B:111:PRO:HD3	1.89	0.53
1:A:195:ILE:CG2	1:A:231:PHE:HA	2.38	0.53
1:C:453:THR:CG2	1:C:471:VAL:HG12	2.38	0.53
1:B:133:ASN:O	1:B:134:TYR:HB2	2.08	0.53
1:B:68:MET:HE3	1:C:64:GLN:HG2	1.89	0.53
1:C:54:SER:HB2	1:C:270:LYS:HE3	1.90	0.53
1:A:233:GLU:HG3	1:A:271:ASP:OD2	2.08	0.53
1:A:100:HIS:CD2	1:A:101:GLN:HG2	2.43	0.53
1:D:133:ASN:O	1:D:134:TYR:HB2	2.08	0.53
1:A:213:MET:HE1	1:A:224:VAL:HG21	1.91	0.53
1:C:110:ILE:HB	1:C:111:PRO:HD3	1.90	0.52
1:B:277:GLY:HA2	1:B:310:LEU:HD21	1.90	0.52
1:D:31:ILE:CD1	4:D:2013:HOH:O	2.57	0.52
1:A:359:ALA:CB	1:A:421:THR:HA	2.39	0.52
1:D:233:GLU:HG3	1:D:271:ASP:OD2	2.09	0.52
1:A:277:GLY:HA2	1:A:310:LEU:HD21	1.90	0.52
1:D:38:PRO:CA	4:D:2013:HOH:O	2.55	0.52
1:D:468:LEU:N	1:D:468:LEU:HD12	2.24	0.52
1:C:100:HIS:CD2	1:C:101:GLN:HG2	2.44	0.52
1:A:53:ASP:OD2	1:A:424:ARG:NH2	2.43	0.52
1:D:344:GLY:HA3	1:D:436:ILE:HD13	1.92	0.52
1:A:344:GLY:HA3	1:A:436:ILE:HD13	1.91	0.51
1:D:213:MET:HE1	1:D:224:VAL:HG21	1.92	0.51
1:D:51:LEU:CD1	1:D:394:VAL:HG13	2.38	0.51
1:D:23:THR:HA	4:D:2010:HOH:O	2.10	0.51
1:C:133:ASN:O	1:C:134:TYR:HB2	2.10	0.51
1:B:53:ASP:OD2	1:B:424:ARG:NH2	2.44	0.51
1:B:264:LEU:C	1:B:264:LEU:HD12	2.31	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:419:ARG:HD3	1:D:421:THR:OG1	2.11	0.51
1:B:359:ALA:CB	1:B:421:THR:HA	2.40	0.51
1:C:233:GLU:HG3	1:C:271:ASP:OD2	2.10	0.51
1:D:359:ALA:CB	1:D:421:THR:HA	2.39	0.51
1:A:468:LEU:N	1:A:468:LEU:HD12	2.26	0.51
1:C:264:LEU:C	1:C:264:LEU:HD12	2.31	0.51
1:D:54:SER:HB2	1:D:270:LYS:HE3	1.93	0.51
1:B:468:LEU:HD12	1:B:468:LEU:N	2.26	0.50
1:D:249:THR:OG1	1:D:252:GLN:HG3	2.12	0.50
1:A:264:LEU:HD12	1:A:264:LEU:C	2.31	0.50
1:C:468:LEU:HD12	1:C:468:LEU:N	2.26	0.50
1:D:344:GLY:O	1:D:347:GLU:CG	2.60	0.50
1:B:334:ARG:HH21	1:B:424:ARG:NH1	2.09	0.50
1:D:469:LYS:HG2	1:D:470:GLU:H	1.76	0.50
1:A:9:GLU:HG3	1:A:333:TYR:CZ	2.47	0.50
1:B:95:TYR:HB3	1:B:291:TYR:CD2	2.47	0.49
1:C:213:MET:HE1	1:C:224:VAL:HG21	1.93	0.49
1:C:359:ALA:CB	1:C:421:THR:HA	2.40	0.49
1:B:233:GLU:HG3	1:B:271:ASP:OD2	2.12	0.49
1:B:346:GLU:HG3	4:B:2104:HOH:O	2.13	0.49
1:B:51:LEU:CD1	1:B:394:VAL:HG13	2.39	0.49
1:C:395:GLU:CD	1:C:399:PHE:HB3	2.32	0.49
1:A:140:GLN:HG3	1:A:150:VAL:HG21	1.95	0.49
1:A:249:THR:OG1	1:A:252:GLN:HG3	2.13	0.49
1:B:395:GLU:CD	1:B:399:PHE:HB3	2.32	0.48
1:D:41:LEU:HD11	4:D:2013:HOH:O	2.13	0.48
1:B:164:ARG:HA	1:B:354:GLN:OE1	2.13	0.48
1:C:461:LEU:HD12	1:D:300:VAL:HG22	1.96	0.48
1:D:9:GLU:HG3	1:D:333:TYR:CZ	2.48	0.48
1:D:311:GLU:HA	4:D:2090:HOH:O	2.13	0.48
1:A:395:GLU:CD	1:A:399:PHE:HB3	2.33	0.48
1:D:419:ARG:HD3	1:D:421:THR:CG2	2.43	0.48
1:D:455:THR:HG21	1:D:469:LYS:HB2	1.95	0.48
1:C:9:GLU:HG3	1:C:333:TYR:CZ	2.49	0.48
1:C:140:GLN:HG3	1:C:150:VAL:HG21	1.96	0.48
1:B:261:ALA:O	1:B:283:LYS:HE3	2.13	0.48
1:D:124:ASP:HA	4:D:2045:HOH:O	2.12	0.48
1:C:51:LEU:CD1	1:C:394:VAL:HG13	2.39	0.48
1:A:196:THR:HG21	4:A:2113:HOH:O	2.14	0.48
1:C:344:GLY:HA3	1:C:436:ILE:HD13	1.95	0.48
1:A:207:LEU:HD11	1:A:257:THR:HA	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:400:LEU:O	1:B:400:LEU:HD23	2.14	0.47
1:B:140:GLN:HG3	1:B:150:VAL:HG21	1.96	0.47
1:C:261:ALA:O	1:C:283:LYS:HE3	2.14	0.47
1:B:344:GLY:HA3	1:B:436:ILE:HD13	1.94	0.47
1:D:261:ALA:O	1:D:283:LYS:HE3	2.13	0.47
1:C:207:LEU:HD11	1:C:257:THR:HA	1.96	0.47
1:A:78:ARG:NH1	4:A:2032:HOH:O	2.47	0.47
1:D:344:GLY:O	1:D:347:GLU:HG3	2.14	0.47
1:B:207:LEU:HD11	1:B:257:THR:HA	1.96	0.47
1:A:419:ARG:HD3	1:A:421:THR:OG1	2.15	0.47
1:C:334:ARG:HH21	1:C:424:ARG:NH1	2.13	0.47
1:D:207:LEU:HD11	1:D:257:THR:HA	1.97	0.47
1:D:140:GLN:HG3	1:D:150:VAL:HG21	1.97	0.47
1:C:95:TYR:HB3	1:C:291:TYR:CD2	2.50	0.47
1:B:211:LYS:HG2	1:B:260:TYR:CZ	2.50	0.47
1:A:334:ARG:HH21	1:A:424:ARG:NH1	2.13	0.47
1:B:213:MET:HE1	1:B:224:VAL:HG21	1.96	0.47
1:A:419:ARG:HD3	1:A:421:THR:CG2	2.45	0.47
1:C:211:LYS:HG2	1:C:260:TYR:CZ	2.49	0.47
1:D:395:GLU:CD	1:D:399:PHE:HB3	2.35	0.46
1:B:431:HIS:CE1	1:D:9:GLU:HB2	2.50	0.46
1:B:9:GLU:HG3	1:B:333:TYR:CZ	2.49	0.46
1:B:419:ARG:HD3	1:B:421:THR:OG1	2.16	0.46
1:D:196:THR:HG23	1:D:202:GLY:HA2	1.98	0.46
1:B:249:THR:OG1	1:B:252:GLN:HG3	2.15	0.46
1:C:419:ARG:HD3	1:C:421:THR:OG1	2.15	0.46
1:A:27:ARG:HG2	1:A:386:TYR:CZ	2.51	0.46
1:C:168:LYS:HE3	1:C:354:GLN:OE1	2.16	0.46
1:C:400:LEU:HD23	1:C:400:LEU:O	2.16	0.46
1:C:249:THR:OG1	1:C:252:GLN:HG3	2.15	0.46
1:D:400:LEU:HD23	1:D:400:LEU:O	2.16	0.46
1:A:261:ALA:O	1:A:283:LYS:HE3	2.14	0.46
1:A:400:LEU:HD23	1:A:400:LEU:O	2.16	0.45
1:A:300:VAL:HG22	1:B:461:LEU:HD12	1.98	0.45
1:A:95:TYR:HB3	1:A:291:TYR:CD2	2.51	0.45
1:D:27:ARG:NH1	1:D:387:LYS:O	2.39	0.45
1:A:195:ILE:HD13	1:A:195:ILE:HA	1.75	0.45
1:A:129:VAL:HG11	1:A:151:ARG:HG3	1.98	0.45
1:A:136:PHE:HB2	1:A:139:THR:OG1	2.17	0.45
1:D:384:GLU:O	1:D:388:VAL:HG23	2.16	0.45
1:D:41:LEU:O	1:D:392:ARG:NH2	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:ASN:O	1:C:255:ARG:NH2	2.50	0.45
1:B:213:MET:HE1	1:B:214:TYR:HA	1.99	0.45
1:C:4:PHE:N	4:C:2001:HOH:O	2.49	0.45
1:A:469:LYS:O	1:A:470:GLU:HB3	2.17	0.45
1:B:288:PHE:O	1:B:292:THR:HG23	2.17	0.45
1:D:288:PHE:O	1:D:292:THR:HG23	2.17	0.45
1:A:27:ARG:NH1	1:A:387:LYS:O	2.48	0.45
1:B:176:LEU:O	1:B:180:ILE:HG13	2.17	0.44
1:A:288:PHE:O	1:A:292:THR:HG23	2.16	0.44
1:A:41:LEU:O	1:A:392:ARG:NH2	2.51	0.44
1:A:395:GLU:O	1:A:396:ILE:HD13	2.17	0.44
1:A:69:ARG:HH21	1:D:69:ARG:NH2	2.08	0.44
1:A:196:THR:HG23	1:A:202:GLY:HA2	2.00	0.44
1:A:164:ARG:HA	1:A:354:GLN:OE1	2.17	0.44
1:A:168:LYS:HE3	1:A:354:GLN:OE1	2.17	0.44
1:D:334:ARG:HH21	1:D:424:ARG:NH1	2.16	0.44
1:B:403:ARG:NH1	1:B:457:GLU:OE2	2.51	0.44
1:D:211:LYS:HG2	1:D:260:TYR:CZ	2.51	0.44
1:B:419:ARG:HD3	1:B:421:THR:CG2	2.44	0.44
1:C:288:PHE:O	1:C:292:THR:HG23	2.18	0.44
1:B:20:LYS:HG2	1:B:21:ARG:N	2.32	0.44
1:C:419:ARG:HD3	1:C:421:THR:CG2	2.46	0.44
1:D:27:ARG:HG2	1:D:386:TYR:CZ	2.53	0.44
1:C:300:VAL:HG22	1:D:461:LEU:HD12	1.99	0.44
1:D:454:PHE:CD2	1:D:457:GLU:HB2	2.53	0.44
1:A:454:PHE:CD2	1:A:457:GLU:HB2	2.53	0.44
1:D:213:MET:HE1	1:D:214:TYR:HA	2.00	0.44
1:C:288:PHE:O	1:C:291:TYR:HB3	2.17	0.44
1:D:95:TYR:HB3	1:D:291:TYR:CD2	2.52	0.44
1:B:196:THR:HG23	1:B:202:GLY:HA2	2.00	0.43
1:C:453:THR:HG22	1:C:471:VAL:HG12	1.99	0.43
1:C:403:ARG:NH1	1:C:457:GLU:OE2	2.51	0.43
1:C:124:ASP:OD1	1:C:126:SER:OG	2.30	0.43
1:C:27:ARG:NH1	1:C:387:LYS:O	2.44	0.43
1:C:384:GLU:O	1:C:388:VAL:HG23	2.18	0.43
1:C:90:ILE:HD13	1:C:326:MET:HE2	2.00	0.43
1:B:168:LYS:HE3	1:B:354:GLN:OE1	2.19	0.43
1:B:41:LEU:O	1:B:392:ARG:NH2	2.52	0.43
1:C:27:ARG:HG2	1:C:386:TYR:CZ	2.54	0.43
1:B:137:ASP:OD2	1:B:138:THR:N	2.52	0.43
1:D:403:ARG:NH1	1:D:457:GLU:OE2	2.52	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:THR:HG23	1:A:63:MET:CE	2.49	0.43
1:A:9:GLU:HB2	1:C:431:HIS:CE1	2.54	0.43
1:A:137:ASP:OD2	1:A:138:THR:N	2.52	0.43
1:A:403:ARG:NH1	1:A:457:GLU:OE2	2.52	0.42
1:B:136:PHE:HB2	1:B:139:THR:OG1	2.18	0.42
1:A:361:PHE:CE2	1:A:419:ARG:HG3	2.55	0.42
1:A:207:LEU:HD23	1:A:245:TYR:CZ	2.54	0.42
1:C:454:PHE:CD2	1:C:457:GLU:HB2	2.54	0.42
1:B:384:GLU:O	1:B:388:VAL:HG23	2.18	0.42
1:B:12:ARG:HH11	1:D:17:GLU:CD	2.23	0.42
1:A:268:ALA:O	1:A:272:ALA:HB3	2.18	0.42
1:C:268:ALA:O	1:C:272:ALA:HB3	2.19	0.42
1:B:194:THR:HA	1:B:227:ASP:HB3	2.01	0.42
1:A:158:ALA:HA	1:A:168:LYS:HB3	2.02	0.42
1:D:176:LEU:O	1:D:180:ILE:HG13	2.19	0.42
1:A:51:LEU:CD1	1:A:394:VAL:HG13	2.38	0.42
1:A:211:LYS:HG2	1:A:260:TYR:CZ	2.53	0.42
1:A:50:LEU:HD11	1:A:435:ILE:HD13	2.02	0.42
1:A:451:GLY:O	1:A:470:GLU:HG2	2.20	0.42
1:B:195:ILE:HD13	1:B:195:ILE:HA	1.76	0.42
1:D:344:GLY:O	1:D:347:GLU:HG2	2.20	0.42
1:C:176:LEU:O	1:C:180:ILE:HG13	2.20	0.42
1:C:213:MET:HE1	1:C:214:TYR:HA	2.02	0.42
1:C:95:TYR:HA	4:C:2038:HOH:O	2.20	0.42
1:D:268:ALA:O	1:D:272:ALA:HB3	2.20	0.42
1:A:469:LYS:O	1:A:471:VAL:N	2.53	0.42
1:A:17:GLU:CD	1:C:12:ARG:HH11	2.23	0.42
1:D:49:ASP:OD2	1:D:392:ARG:CD	2.68	0.41
1:B:361:PHE:CE2	1:B:419:ARG:HG3	2.55	0.41
1:D:168:LYS:HE3	1:D:354:GLN:OE1	2.20	0.41
1:B:361:PHE:CZ	1:B:419:ARG:HG3	2.55	0.41
1:A:194:THR:HA	1:A:227:ASP:HB3	2.01	0.41
1:C:196:THR:HG23	1:C:202:GLY:HA2	2.02	0.41
1:C:164:ARG:HA	1:C:354:GLN:OE1	2.20	0.41
1:D:469:LYS:HG2	1:D:470:GLU:N	2.35	0.41
1:D:469:LYS:O	1:D:471:VAL:N	2.53	0.41
1:A:361:PHE:CZ	1:A:419:ARG:HG3	2.56	0.41
1:D:194:THR:HA	1:D:227:ASP:HB3	2.02	0.41
1:C:120:GLU:CB	4:C:2041:HOH:O	2.69	0.41
1:C:41:LEU:O	1:C:392:ARG:NH2	2.53	0.41
1:C:207:LEU:HD23	1:C:245:TYR:CZ	2.56	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:207:LEU:HD23	1:D:245:TYR:CZ	2.55	0.41
1:B:14:ARG:NE	1:B:14:ARG:HA	2.35	0.41
1:A:49:ASP:OD2	1:A:392:ARG:CD	2.69	0.41
1:A:452:LEU:CD2	1:A:471:VAL:N	2.76	0.41
1:C:60:THR:OG1	1:C:63:MET:HG3	2.21	0.41
1:C:361:PHE:CZ	1:C:419:ARG:HG3	2.56	0.41
1:C:361:PHE:CE2	1:C:419:ARG:HG3	2.56	0.41
1:B:288:PHE:O	1:B:291:TYR:HB3	2.20	0.41
1:B:158:ALA:HA	1:B:168:LYS:HB3	2.02	0.41
1:C:211:LYS:HE2	1:C:260:TYR:OH	2.21	0.41
1:B:454:PHE:CD2	1:B:457:GLU:HB2	2.56	0.41
1:B:27:ARG:HG2	1:B:386:TYR:CZ	2.56	0.41
1:D:288:PHE:O	1:D:291:TYR:HB3	2.21	0.41
1:D:117:ARG:HG3	1:D:117:ARG:HH11	1.86	0.41
1:D:14:ARG:NE	1:D:14:ARG:HA	2.36	0.41
1:B:108:ILE:O	1:B:111:PRO:HD2	2.22	0.40
1:C:108:ILE:O	1:C:111:PRO:HD2	2.21	0.40
1:B:195:ILE:HG23	1:B:234:ASN:CB	2.52	0.40
1:B:211:LYS:HE2	1:B:260:TYR:OH	2.22	0.40
1:A:384:GLU:O	1:A:388:VAL:HG23	2.20	0.40
1:A:176:LEU:O	1:A:180:ILE:HG13	2.20	0.40
1:C:194:THR:HA	1:C:227:ASP:HB3	2.03	0.40
1:D:184:GLY:HA2	1:D:185:PRO:HD2	1.95	0.40
4:A:2013:HOH:O	1:B:75:SER:HB3	2.21	0.40
1:C:125:ARG:HG2	1:C:128:MET:HE1	2.04	0.40
1:C:195:ILE:HA	1:C:195:ILE:HD13	1.76	0.40
1:B:226:MET:HG2	1:B:258:TYR:CE1	2.57	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	465/471 (99%)	441 (95%)	22 (5%)	2 (0%)	39	74
1	B	465/471 (99%)	446 (96%)	17 (4%)	2 (0%)	39	74
1	C	466/471 (99%)	445 (96%)	19 (4%)	2 (0%)	39	74
1	D	465/471 (99%)	442 (95%)	21 (4%)	2 (0%)	39	74
All	All	1861/1884 (99%)	1774 (95%)	79 (4%)	8 (0%)	39	74

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	134	TYR
1	B	134	TYR
1	C	134	TYR
1	D	134	TYR
1	B	196	THR
1	C	196	THR
1	A	196	THR
1	D	196	THR

### 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	360/391 (92%)	348 (97%)	12 (3%)	45	79
1	B	362/391 (93%)	349 (96%)	13 (4%)	42	76
1	C	361/391 (92%)	349 (97%)	12 (3%)	45	79
1	D	352/391 (90%)	337 (96%)	15 (4%)	35	70
All	All	1435/1564 (92%)	1383 (96%)	52 (4%)	42	76

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

Mol	Chain	Res	Type
1	A	14	ARG
1	A	69	ARG
1	A	134	TYR

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Mol	Chain	Res	Type
1	A	196	THR
1	A	213	MET
1	A	292	THR
1	A	334	ARG
1	A	379	GLN
1	A	392	ARG
1	A	394	VAL
1	A	395	GLU
1	A	421	THR
1	B	14	ARG
1	B	24	ARG
1	B	134	TYR
1	B	196	THR
1	B	213	MET
1	B	292	THR
1	B	334	ARG
1	B	379	GLN
1	B	392	ARG
1	B	394	VAL
1	B	395	GLU
1	B	421	THR
1	B	470	GLU
1	C	14	ARG
1	C	24	ARG
1	C	134	TYR
1	C	196	THR
1	C	213	MET
1	C	292	THR
1	C	334	ARG
1	C	379	GLN
1	C	392	ARG
1	C	394	VAL
1	C	395	GLU
1	C	421	THR
1	D	14	ARG
1	D	24	ARG
1	D	103	ARG
1	D	134	TYR
1	D	196	THR
1	D	213	MET
1	D	292	THR
1	D	334	ARG

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Mol	Chain	Res	Type
1	D	379	GLN
1	D	392	ARG
1	D	394	VAL
1	D	395	GLU
1	D	421	THR
1	D	462	ARG
1	D	470	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	339	GLN
1	A	379	GLN
1	B	379	GLN
1	C	379	GLN
1	D	379	GLN

### 5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

## 5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 6 are monoatomic - leaving 8 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	SO4	A	500	-	4,4,4	1.69	0	6,6,6	1.57	1 (16%)
2	SO4	A	501	-	4,4,4	1.72	1 (25%)	6,6,6	1.62	1 (16%)
2	SO4	B	500	-	4,4,4	1.59	0	6,6,6	1.60	1 (16%)
2	SO4	B	501	-	4,4,4	1.64	1 (25%)	6,6,6	1.62	1 (16%)
2	SO4	C	500	-	4,4,4	1.73	0	6,6,6	1.56	1 (16%)
2	SO4	C	501	-	4,4,4	1.55	0	6,6,6	1.61	1 (16%)
2	SO4	D	500	-	4,4,4	1.66	1 (25%)	6,6,6	1.64	1 (16%)
2	SO4	D	501	-	4,4,4	1.53	0	6,6,6	1.60	1 (16%)

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	500	-	-	0/0/0/0	0/0/0/0
2	SO4	A	501	-	-	0/0/0/0	0/0/0/0
2	SO4	B	500	-	-	0/0/0/0	0/0/0/0
2	SO4	B	501	-	-	0/0/0/0	0/0/0/0
2	SO4	C	500	-	-	0/0/0/0	0/0/0/0
2	SO4	C	501	-	-	0/0/0/0	0/0/0/0
2	SO4	D	500	-	-	0/0/0/0	0/0/0/0
2	SO4	D	501	-	-	0/0/0/0	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	SO4	O2-S	2.03	1.54	1.47
2	D	500	SO4	O1-S	2.09	1.54	1.47
2	A	501	SO4	O2-S	2.14	1.54	1.47

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	C	500	SO4	O4-S-O3	3.64	123.80	108.98
2	A	500	SO4	O4-S-O3	3.69	124.00	108.98
2	D	501	SO4	O4-S-O3	3.72	124.11	108.98
2	C	501	SO4	O4-S-O3	3.72	124.12	108.98
2	A	501	SO4	O4-S-O3	3.76	124.28	108.98
2	B	500	SO4	O4-S-O3	3.77	124.31	108.98

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	SO4	O4-S-O3	3.77	124.32	108.98
2	D	500	SO4	O4-S-O3	3.85	124.63	108.98

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	500	SO4	1	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, 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	467/471 (99%)	-0.49	5 (1%) 82 74	6, 12, 24, 35	0
1	B	467/471 (99%)	-0.40	2 (0%) 93 90	6, 12, 24, 34	0
1	C	468/471 (99%)	-0.44	0 100 100	6, 13, 24, 35	0
1	D	467/471 (99%)	-0.08	32 (6%) 20 11	6, 13, 25, 36	0
All	All	1869/1884 (99%)	-0.35	39 (2%) 67 56	6, 12, 24, 36	0

All (39) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	471	VAL	4.9
1	D	405	PRO	4.6
1	D	404	ASP	4.4
1	D	406	LYS	4.0
1	D	407	THR	3.9
1	D	411	LEU	3.5
1	D	412	PRO	3.3
1	D	372	PRO	3.3
1	B	471	VAL	3.2
1	D	413	CYS	3.1
1	D	446	ALA	3.1
1	D	447	SER	3.0
1	D	453	THR	3.0
1	D	445	ASN	2.9
1	D	409	LYS	2.9
1	D	449	ILE	2.9
1	D	369	PRO	2.9
1	D	443	LYS	2.8
1	D	469	LYS	2.8
1	D	410	GLN	2.7
1	D	441	HIS	2.6

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Mol	Chain	Res	Type	RSRZ
1	D	374	ASP	2.5
1	B	243	ALA	2.5
1	D	347	GLU	2.4
1	D	365	GLY	2.4
1	A	405	PRO	2.3
1	A	408	GLY	2.3
1	A	412	PRO	2.3
1	D	370	HIS	2.3
1	A	407	THR	2.2
1	D	448	ASN	2.2
1	D	26	TYR	2.2
1	D	373	ALA	2.2
1	D	402	GLY	2.1
1	D	455	THR	2.1
1	D	35	GLY	2.1
1	D	25	ALA	2.0
1	D	349	GLY	2.0
1	A	406	LYS	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 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	RSCC	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	SO4	A	501	5/5	0.97	0.14	-0.44	23,24,26,33	0
2	SO4	D	501	5/5	0.96	0.15	-0.46	50,51,53,56	0
3	K	B	1472	1/1	0.96	0.13	-0.57	27,27,27,27	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	SO4	D	500	5/5	0.99	0.13	-1.15	11,16,19,20	0
2	SO4	B	501	5/5	0.98	0.11	-1.40	25,28,30,32	0
2	SO4	C	501	5/5	0.98	0.12	-2.02	22,25,30,31	0
3	K	D	1471	1/1	0.98	0.11	-2.08	14,14,14,14	0
2	SO4	A	500	5/5	0.98	0.10	-2.12	11,11,15,15	0
2	SO4	C	500	5/5	0.99	0.11	-2.17	11,14,16,17	0
3	K	A	1472	1/1	0.97	0.10	-2.56	19,19,19,19	0
2	SO4	B	500	5/5	0.98	0.09	-2.89	22,24,25,26	0
3	K	A	1471	1/1	0.96	0.11	-3.12	18,18,18,18	0
3	K	C	1472	1/1	0.95	0.09	-4.49	13,13,13,13	0
3	K	C	1473	1/1	0.96	0.06	-6.87	19,19,19,19	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.