



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 31, 2021 – 11:34 PM EDT

PDB ID : 2DTZ
Title : Crystal Structure of multidrug binding protein QacR from *Staphylococcus aureus* cocrystallized with compound DB75
Authors : Brooks, B.E.; Brennan, R.G.
Deposited on : 2006-07-19
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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.23.2
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

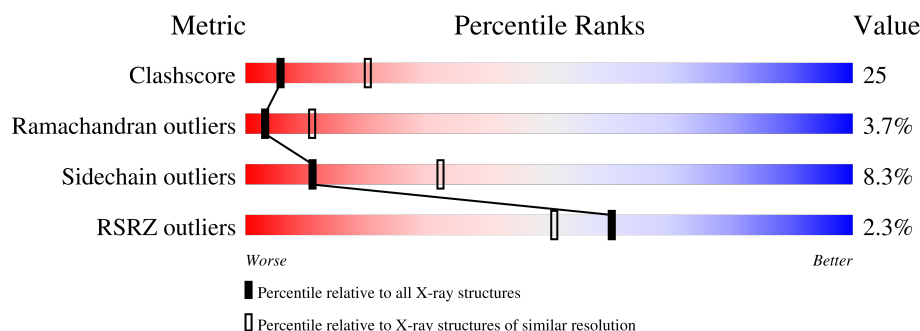
1 Overall quality at a glance

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))
Clashscore	141614	3569 (2.80-2.80)
Ramachandran outliers	138981	3498 (2.80-2.80)
Sidechain outliers	138945	3500 (2.80-2.80)
RSRZ outliers	127900	3078 (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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	194	
1	B	194	
1	D	194	
1	E	194	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6063 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 HTH-type transcriptional regulator qacR.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	B	182	Total	C	N	O	S	0	0	0
			1477	950	241	284	2			
1	D	184	Total	C	N	O	S	0	0	0
			1466	945	240	279	2			
1	A	186	Total	C	N	O	S	0	0	0
			1529	986	250	291	2			
1	E	186	Total	C	N	O	S	0	0	0
			1528	985	248	293	2			

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	72	ALA	CYS	engineered mutation	UNP P0A0N4
B	141	SER	CYS	engineered mutation	UNP P0A0N4
B	189	HIS	-	expression tag	UNP P0A0N4
B	190	HIS	-	expression tag	UNP P0A0N4
B	191	HIS	-	expression tag	UNP P0A0N4
B	192	HIS	-	expression tag	UNP P0A0N4
B	193	HIS	-	expression tag	UNP P0A0N4
B	194	HIS	-	expression tag	UNP P0A0N4
D	72	ALA	CYS	engineered mutation	UNP P0A0N4
D	141	SER	CYS	engineered mutation	UNP P0A0N4
D	189	HIS	-	expression tag	UNP P0A0N4
D	190	HIS	-	expression tag	UNP P0A0N4
D	191	HIS	-	expression tag	UNP P0A0N4
D	192	HIS	-	expression tag	UNP P0A0N4
D	193	HIS	-	expression tag	UNP P0A0N4
D	194	HIS	-	expression tag	UNP P0A0N4
A	72	ALA	CYS	engineered mutation	UNP P0A0N4
A	141	SER	CYS	engineered mutation	UNP P0A0N4
A	189	HIS	-	expression tag	UNP P0A0N4
A	190	HIS	-	expression tag	UNP P0A0N4
A	191	HIS	-	expression tag	UNP P0A0N4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	192	HIS	-	expression tag	UNP P0A0N4
A	193	HIS	-	expression tag	UNP P0A0N4
A	194	HIS	-	expression tag	UNP P0A0N4
E	72	ALA	CYS	engineered mutation	UNP P0A0N4
E	141	SER	CYS	engineered mutation	UNP P0A0N4
E	189	HIS	-	expression tag	UNP P0A0N4
E	190	HIS	-	expression tag	UNP P0A0N4
E	191	HIS	-	expression tag	UNP P0A0N4
E	192	HIS	-	expression tag	UNP P0A0N4
E	193	HIS	-	expression tag	UNP P0A0N4
E	194	HIS	-	expression tag	UNP P0A0N4

- Molecule 2 is SULFATE ION (three-letter code: SO₄) (formula: O₄S).



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

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

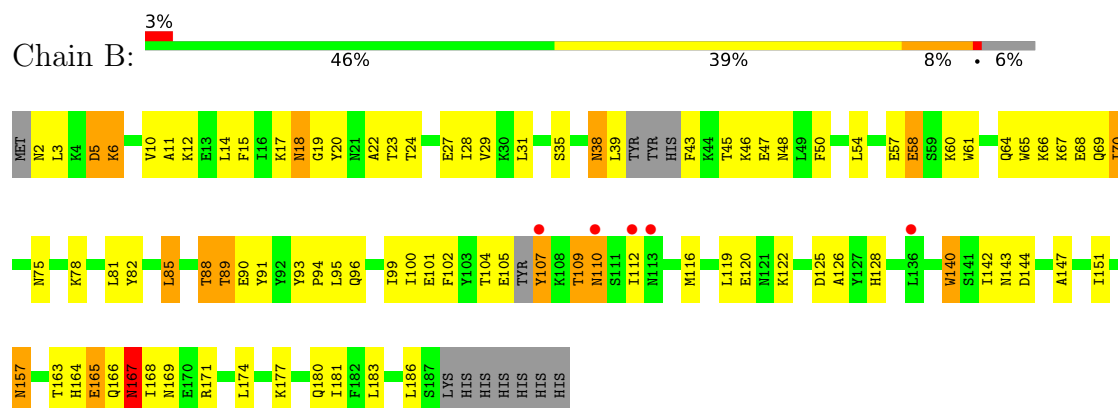
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	2	Total	O	0	0
			2	2		
3	A	9	Total	O	0	0
			9	9		
3	E	12	Total	O	0	0
			12	12		

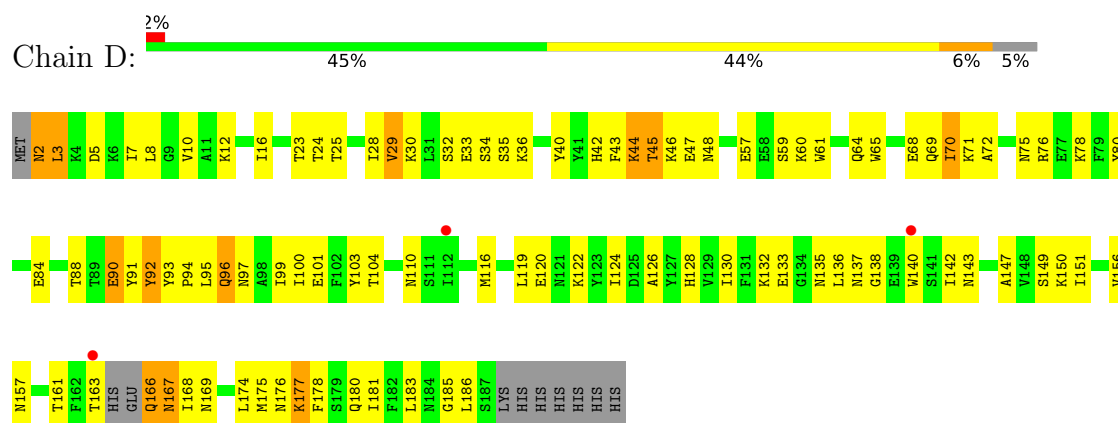
3 Residue-property plots

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

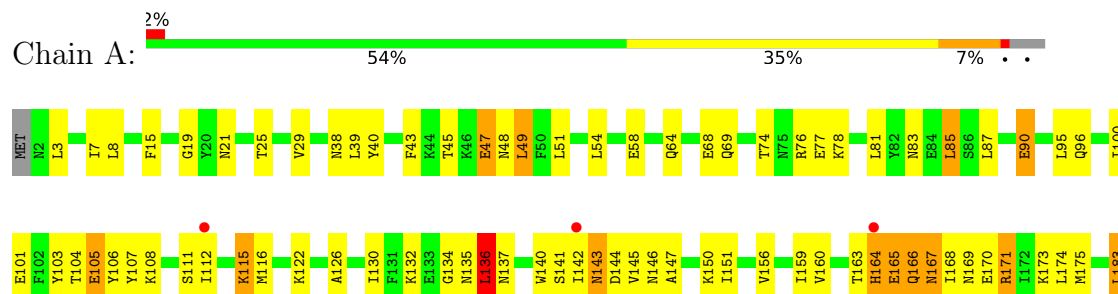
- Molecule 1: HTH-type transcriptional regulator qacR



- Molecule 1: HTH-type transcriptional regulator qacR

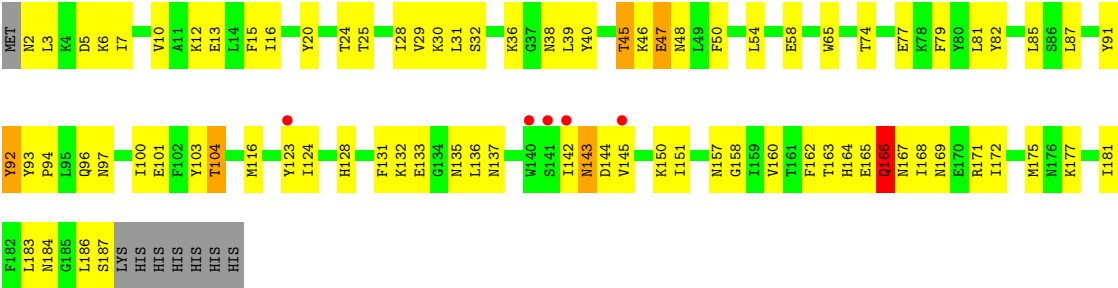


- Molecule 1: HTH-type transcriptional regulator qacR





● Molecule 1: HTH-type transcriptional regulator qacR



4 Data and refinement statistics

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α , β , γ	171.50Å 171.50Å 94.32Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	76.70 – 2.80 76.70 – 2.80	Depositor EDS
% Data completeness (in resolution range)	97.9 (76.70-2.80) 98.0 (76.70-2.80)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.88 (at 2.82Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.224 , 0.289 0.223 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	69.1	Xtriage
Anisotropy	0.249	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 75.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6063	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.08% 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: 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	0.45	0/1559	0.64	0/2102
1	B	0.40	0/1502	0.58	0/2025
1	D	0.40	0/1494	0.58	0/2024
1	E	0.43	0/1557	0.62	0/2099
All	All	0.42	0/6112	0.61	0/8250

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1529	0	1510	75	0
1	B	1477	0	1445	86	0
1	D	1466	0	1397	78	0
1	E	1528	0	1509	79	0
2	A	20	0	0	0	0
2	D	5	0	0	0	0
2	E	15	0	0	0	0
3	A	9	0	0	0	0
3	D	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	E	12	0	0	0	0
All	All	6063	0	5861	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 25.

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:A:163:THR:O	1:A:165:GLU:HG2	1.61	1.00
1:A:135:ASN:HD21	1:A:142:ILE:H	1.00	0.97
1:B:96:GLN:HG3	1:B:157:ASN:HD21	1.29	0.96
1:A:68:GLU:HG2	1:A:85:LEU:HD21	1.50	0.93
1:B:39:LEU:H	1:B:39:LEU:HD22	1.36	0.89
1:D:126:ALA:O	1:D:130:ILE:HG12	1.75	0.87
1:B:164:HIS:HA	1:B:171:ARG:NH1	1.90	0.86
1:A:115:LYS:HD2	1:A:115:LYS:H	1.40	0.86
1:D:101:GLU:HG3	1:E:97:ASN:HD21	1.42	0.85
1:A:143:ASN:HD22	1:A:143:ASN:H	1.24	0.85
1:D:45:THR:HG23	1:D:48:ASN:HB2	1.59	0.84
1:A:76:ARG:HG3	1:A:183:LEU:HD23	1.61	0.83
1:E:3:LEU:O	1:E:7:ILE:HG12	1.81	0.80
1:E:45:THR:HG23	1:E:48:ASN:HB2	1.65	0.78
1:B:177:LYS:HE2	1:A:144:ASP:OD2	1.83	0.78
1:A:135:ASN:HD21	1:A:142:ILE:N	1.82	0.76
1:A:83:ASN:ND2	1:A:156:VAL:HG21	2.01	0.75
1:D:5:ASP:HA	1:D:8:LEU:HD12	1.68	0.75
1:A:142:ILE:HD11	1:A:186:LEU:HD13	1.69	0.75
1:B:6:LYS:HE3	1:B:6:LYS:HA	1.67	0.75
1:B:54:LEU:HB3	1:B:119:LEU:HD11	1.67	0.75
1:D:147:ALA:O	1:D:151:ILE:HG13	1.87	0.75
1:E:143:ASN:ND2	1:E:144:ASP:H	1.85	0.74
1:A:146:ASN:O	1:A:150:LYS:HG3	1.88	0.73
1:D:96:GLN:O	1:D:100:ILE:HG12	1.89	0.73
1:D:163:THR:HG22	1:D:166:GLN:HG3	1.71	0.73
1:B:89:THR:HG22	1:B:90:GLU:H	1.53	0.72
1:A:164:HIS:HA	1:A:171:ARG:NH1	2.04	0.72
1:E:87:LEU:HD23	1:E:160:VAL:HG22	1.70	0.72
1:D:76:ARG:HG3	1:D:183:LEU:HD13	1.70	0.72
1:E:165:GLU:HB2	1:E:166:GLN:NE2	2.06	0.71
1:B:45:THR:OG1	1:B:48:ASN:HB2	1.91	0.71

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:65:TRP:O	1:D:69:GLN:HB3	1.91	0.70
1:E:142:ILE:HD11	1:E:186:LEU:HD13	1.73	0.70
1:A:143:ASN:H	1:A:143:ASN:ND2	1.89	0.70
1:E:143:ASN:CG	1:E:144:ASP:H	1.96	0.69
1:E:74:THR:OG1	1:E:77:GLU:HG3	1.92	0.69
1:B:95:LEU:O	1:B:99:ILE:HG12	1.94	0.68
1:B:66:LYS:C	1:B:68:GLU:H	1.98	0.67
1:E:100:ILE:O	1:E:104:THR:HG23	1.94	0.66
1:D:24:THR:O	1:D:28:ILE:HG12	1.97	0.65
1:A:135:ASN:ND2	1:A:142:ILE:H	1.84	0.65
1:B:75:ASN:HA	1:B:78:LYS:HG3	1.79	0.65
1:D:157:ASN:O	1:D:161:THR:HG23	1.97	0.65
1:E:45:THR:OG1	1:E:47:GLU:HG2	1.97	0.65
1:A:147:ALA:O	1:A:151:ILE:HG12	1.97	0.65
1:A:134:GLY:HA3	1:A:140:TRP:CE2	2.32	0.64
1:A:69:GLN:HB2	1:A:78:LYS:HD3	1.79	0.64
1:E:128:HIS:HD2	1:E:145:VAL:HG11	1.63	0.63
1:B:177:LYS:O	1:B:181:ILE:HG12	1.99	0.63
1:A:136:LEU:N	1:A:136:LEU:CD1	2.62	0.63
1:E:142:ILE:CD1	1:E:186:LEU:HD13	2.27	0.63
1:D:100:ILE:HD11	1:E:162:PHE:CE2	2.35	0.62
1:E:20:TYR:CZ	1:E:46:LYS:HE2	2.35	0.62
1:D:177:LYS:HD2	1:D:181:ILE:HD11	1.81	0.61
1:D:116:MET:O	1:D:120:GLU:HG2	2.00	0.61
1:A:126:ALA:O	1:A:130:ILE:HG12	2.01	0.61
1:E:163:THR:O	1:E:171:ARG:HD3	2.01	0.61
1:D:57:GLU:HG3	1:D:95:LEU:HD12	1.83	0.60
1:D:45:THR:CG2	1:D:48:ASN:HB2	2.31	0.60
1:E:97:ASN:O	1:E:101:GLU:HG3	2.01	0.60
1:D:177:LYS:O	1:D:177:LYS:HD3	2.01	0.60
1:D:7:ILE:HD13	1:D:32:SER:OG	2.01	0.60
1:E:163:THR:C	1:E:165:GLU:H	2.05	0.60
1:A:136:LEU:N	1:A:136:LEU:HD12	2.16	0.60
1:E:166:GLN:NE2	1:E:166:GLN:H	1.98	0.59
1:A:165:GLU:O	1:A:166:GLN:HB3	2.01	0.59
1:D:59:SER:C	1:D:61:TRP:H	2.06	0.59
1:B:39:LEU:H	1:B:39:LEU:CD2	2.13	0.59
1:A:164:HIS:O	1:A:165:GLU:HB3	2.02	0.59
1:D:70:ILE:C	1:D:72:ALA:H	2.05	0.58
1:D:95:LEU:O	1:D:99:ILE:HG13	2.04	0.58
1:D:2:ASN:ND2	1:D:3:LEU:N	2.52	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:58:GLU:HG2	1:A:122:LYS:HD2	1.86	0.58
1:B:64:GLN:O	1:B:68:GLU:HB2	2.03	0.58
1:D:88:THR:HG22	1:D:168:ILE:HD11	1.85	0.58
1:E:163:THR:HA	1:E:165:GLU:OE1	2.04	0.58
1:E:92:TYR:CD2	1:E:123:TYR:CZ	2.93	0.57
1:A:87:LEU:O	1:A:87:LEU:HD12	2.04	0.57
1:A:136:LEU:CD1	1:A:136:LEU:H	2.17	0.57
1:B:14:LEU:HD22	1:B:28:ILE:HD13	1.86	0.57
1:B:89:THR:HG22	1:B:90:GLU:N	2.20	0.57
1:B:116:MET:O	1:B:120:GLU:HG3	2.04	0.56
1:B:167:ASN:ND2	1:B:169:ASN:H	2.03	0.56
1:B:24:THR:CG2	1:B:27:GLU:HG3	2.36	0.56
1:D:84:GLU:CG	1:D:176:ASN:HD21	2.18	0.56
1:D:177:LYS:O	1:D:181:ILE:HG13	2.06	0.56
1:E:45:THR:HG23	1:E:48:ASN:CB	2.33	0.56
1:D:80:TYR:OH	1:D:180:GLN:HB2	2.06	0.56
1:D:156:VAL:HG13	1:D:175:MET:CE	2.35	0.56
1:A:51:LEU:HD11	1:A:112:ILE:HG23	1.87	0.56
1:E:79:PHE:CD2	1:E:183:LEU:HD13	2.40	0.56
1:B:96:GLN:NE2	1:B:157:ASN:OD1	2.38	0.55
1:B:167:ASN:HD22	1:B:167:ASN:C	2.07	0.55
1:A:164:HIS:HA	1:A:171:ARG:CZ	2.37	0.55
1:A:106:TYR:CD1	1:A:112:ILE:HG13	2.42	0.55
1:A:135:ASN:C	1:A:137:ASN:H	2.10	0.55
1:E:96:GLN:NE2	1:E:157:ASN:HD21	2.05	0.55
1:A:134:GLY:HA3	1:A:140:TRP:CZ2	2.42	0.54
1:B:47:GLU:OE2	1:B:112:ILE:HD11	2.08	0.54
1:B:15:PHE:CE1	1:B:50:PHE:HD2	2.24	0.54
1:B:109:THR:HG23	1:B:112:ILE:HG12	1.90	0.54
1:D:136:LEU:C	1:D:138:GLY:H	2.11	0.54
1:D:142:ILE:HD11	1:D:186:LEU:HD13	1.89	0.54
1:B:66:LYS:HD3	1:B:69:GLN:HE21	1.71	0.54
1:B:125:ASP:O	1:B:128:HIS:N	2.40	0.54
1:D:119:LEU:O	1:D:122:LYS:N	2.41	0.54
1:A:135:ASN:O	1:A:137:ASN:N	2.41	0.54
1:B:89:THR:HG22	1:B:91:TYR:H	1.73	0.54
1:A:21:ASN:ND2	1:A:101:GLU:CG	2.70	0.54
1:D:2:ASN:CG	1:D:3:LEU:H	2.11	0.53
1:E:3:LEU:HD13	1:E:38:ASN:ND2	2.24	0.53
1:B:54:LEU:O	1:B:58:GLU:HB2	2.09	0.52
1:A:165:GLU:O	1:A:166:GLN:CB	2.58	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:143:ASN:ND2	1:E:144:ASP:N	2.57	0.52
1:A:21:ASN:ND2	1:A:101:GLU:HG2	2.25	0.52
1:D:151:ILE:HD11	1:E:177:LYS:HB3	1.91	0.52
1:D:151:ILE:CD1	1:E:177:LYS:HB3	2.40	0.52
1:B:119:LEU:HD23	1:B:119:LEU:O	2.10	0.51
1:B:109:THR:HG23	1:B:112:ILE:CG1	2.40	0.51
1:A:15:PHE:O	1:A:19:GLY:N	2.43	0.51
1:B:31:LEU:O	1:B:31:LEU:HG	2.10	0.50
1:B:61:TRP:CZ2	1:B:82:TYR:CE1	2.99	0.50
1:A:143:ASN:ND2	1:A:143:ASN:N	2.53	0.50
1:A:167:ASN:OD1	1:A:170:GLU:HB2	2.11	0.50
1:E:97:ASN:O	1:E:100:ILE:HG22	2.11	0.50
1:D:174:LEU:HB3	1:E:151:ILE:CD1	2.41	0.50
1:A:135:ASN:ND2	1:A:141:SER:HA	2.27	0.50
1:E:96:GLN:NE2	1:E:157:ASN:ND2	2.60	0.50
1:A:165:GLU:O	1:A:165:GLU:HG3	2.11	0.50
1:E:128:HIS:CD2	1:E:145:VAL:HG11	2.45	0.50
1:D:96:GLN:HG2	1:D:161:THR:HG21	1.93	0.50
1:D:96:GLN:NE2	1:D:157:ASN:HD21	2.10	0.50
1:B:88:THR:O	1:B:89:THR:O	2.30	0.50
1:B:66:LYS:C	1:B:68:GLU:N	2.64	0.50
1:E:132:LYS:O	1:E:135:ASN:N	2.45	0.50
1:D:175:MET:O	1:D:178:PHE:HB3	2.11	0.49
1:A:38:ASN:ND2	1:E:13:GLU:CD	2.66	0.49
1:B:90:GLU:N	1:B:90:GLU:OE1	2.46	0.49
1:D:101:GLU:HG3	1:E:97:ASN:ND2	2.21	0.49
1:B:104:THR:C	1:B:105:GLU:HG3	2.32	0.49
1:B:107:TYR:C	1:B:107:TYR:CD2	2.85	0.49
1:D:167:ASN:ND2	1:D:169:ASN:H	2.09	0.49
1:D:97:ASN:HD21	1:E:100:ILE:HG23	1.77	0.49
1:E:143:ASN:CG	1:E:144:ASP:N	2.65	0.49
1:B:20:TYR:N	1:B:101:GLU:OE1	2.46	0.49
1:D:92:TYR:O	1:D:94:PRO:N	2.46	0.49
1:D:10:VAL:HG21	1:D:32:SER:HB3	1.94	0.49
1:D:128:HIS:HA	1:D:149:SER:OG	2.12	0.49
1:A:143:ASN:HD22	1:A:143:ASN:N	1.90	0.49
1:A:115:LYS:HD2	1:A:115:LYS:N	2.18	0.48
1:A:103:TYR:O	1:A:105:GLU:N	2.39	0.48
1:E:30:LYS:C	1:E:32:SER:H	2.16	0.48
1:D:91:TYR:O	1:D:92:TYR:HB3	2.14	0.48
1:D:133:GLU:OE2	1:D:136:LEU:HD12	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:183:LEU:C	1:D:185:GLY:H	2.17	0.48
1:A:103:TYR:C	1:A:105:GLU:H	2.16	0.48
1:E:6:LYS:O	1:E:10:VAL:HG23	2.13	0.48
1:E:58:GLU:HG2	1:E:123:TYR:CE2	2.49	0.48
1:B:6:LYS:HE3	1:B:6:LYS:CA	2.42	0.48
1:D:57:GLU:HG3	1:D:95:LEU:CD1	2.42	0.48
1:A:134:GLY:HA3	1:A:140:TRP:NE1	2.28	0.48
1:A:163:THR:O	1:A:164:HIS:C	2.52	0.48
1:A:168:ILE:HA	1:A:171:ARG:HG3	1.95	0.48
1:E:96:GLN:CD	1:E:157:ASN:HD21	2.16	0.48
1:A:39:LEU:O	1:A:39:LEU:HD23	2.14	0.47
1:B:81:LEU:HG	1:B:85:LEU:CD2	2.44	0.47
1:B:93:TYR:N	1:B:94:PRO:CD	2.76	0.47
1:E:29:VAL:HG23	1:E:39:LEU:HD12	1.96	0.47
1:B:3:LEU:HD13	1:B:3:LEU:O	2.13	0.47
1:B:174:LEU:HB3	1:A:151:ILE:CD1	2.44	0.47
1:A:43:PHE:O	1:A:48:ASN:HB3	2.15	0.47
1:E:10:VAL:HG11	1:E:31:LEU:HB2	1.96	0.47
1:E:186:LEU:O	1:E:187:SER:C	2.53	0.47
1:A:21:ASN:ND2	1:A:101:GLU:OE1	2.48	0.47
1:B:10:VAL:O	1:B:14:LEU:HD13	2.14	0.47
1:A:136:LEU:H	1:A:136:LEU:HD13	1.79	0.47
1:B:142:ILE:HD12	1:B:186:LEU:HD22	1.96	0.47
1:B:167:ASN:ND2	1:B:167:ASN:C	2.68	0.47
1:D:177:LYS:HD3	1:D:177:LYS:C	2.34	0.47
1:E:103:TYR:HD2	1:E:116:MET:HG2	1.80	0.47
1:D:174:LEU:HB3	1:E:151:ILE:HD11	1.97	0.47
1:B:2:ASN:O	1:B:6:LYS:HD2	2.15	0.46
1:B:70:ILE:H	1:B:70:ILE:HG13	1.40	0.46
1:A:25:THR:O	1:A:29:VAL:HG23	2.15	0.46
1:E:25:THR:HA	1:E:28:ILE:HD12	1.96	0.46
1:B:65:TRP:O	1:B:69:GLN:N	2.46	0.46
1:D:12:LYS:O	1:D:16:ILE:HG13	2.15	0.46
1:B:167:ASN:HD22	1:B:168:ILE:N	2.14	0.46
1:D:166:GLN:HE21	1:D:166:GLN:CA	2.29	0.46
1:B:66:LYS:O	1:B:68:GLU:N	2.49	0.46
1:B:125:ASP:O	1:B:126:ALA:C	2.54	0.46
1:A:186:LEU:O	1:A:187:SER:HB3	2.16	0.46
1:E:45:THR:OG1	1:E:46:LYS:N	2.49	0.46
1:B:6:LYS:HA	1:B:6:LYS:CE	2.40	0.46
1:A:163:THR:O	1:A:165:GLU:N	2.49	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:120:GLU:O	1:D:124:ILE:HG13	2.15	0.46
1:D:101:GLU:OE2	1:E:101:GLU:OE2	2.35	0.45
1:A:8:LEU:HD23	1:A:49:LEU:HD21	1.97	0.45
1:B:20:TYR:CE1	1:B:46:LYS:HE2	2.50	0.45
1:B:61:TRP:HZ2	1:B:82:TYR:CE1	2.34	0.45
1:B:38:ASN:O	1:B:39:LEU:C	2.54	0.45
1:A:21:ASN:HB2	1:A:101:GLU:OE1	2.17	0.45
1:D:23:THR:CG2	1:D:28:ILE:HD11	2.46	0.45
1:E:81:LEU:HG	1:E:85:LEU:HD21	1.99	0.45
1:B:24:THR:HG22	1:B:27:GLU:OE2	2.17	0.45
1:E:158:GLY:O	1:E:162:PHE:HD1	1.99	0.45
1:B:147:ALA:O	1:B:151:ILE:HG12	2.16	0.45
1:B:165:GLU:O	1:B:166:GLN:HB3	2.16	0.45
1:D:103:TYR:OH	1:E:165:GLU:HG3	2.17	0.45
1:B:110:ASN:C	1:B:112:ILE:H	2.20	0.44
1:D:64:GLN:O	1:D:68:GLU:HG3	2.17	0.44
1:D:2:ASN:CG	1:D:3:LEU:N	2.71	0.44
1:B:39:LEU:HB3	1:B:43:PHE:O	2.17	0.44
1:A:3:LEU:O	1:A:7:ILE:HG13	2.17	0.44
1:E:12:LYS:O	1:E:16:ILE:HG13	2.17	0.44
1:B:29:VAL:CG1	1:B:35:SER:HA	2.47	0.44
1:B:19:GLY:HA3	1:B:101:GLU:OE1	2.18	0.44
1:A:40:TYR:CZ	1:E:6:LYS:HD3	2.52	0.44
1:E:25:THR:HG21	1:E:40:TYR:OH	2.18	0.44
1:B:11:ALA:HA	1:B:28:ILE:CD1	2.48	0.44
1:B:18:ASN:HB2	1:B:22:ALA:HB3	2.00	0.44
1:B:60:LYS:HD3	1:B:91:TYR:CZ	2.52	0.44
1:E:91:TYR:O	1:E:93:TYR:N	2.47	0.44
1:B:99:ILE:O	1:B:102:PHE:HB3	2.18	0.43
1:D:70:ILE:C	1:D:72:ALA:N	2.71	0.43
1:B:47:GLU:CD	1:B:112:ILE:HD11	2.39	0.43
1:D:25:THR:CG2	1:D:46:LYS:HB2	2.48	0.43
1:D:124:ILE:HD13	1:D:150:LYS:HG3	2.00	0.43
1:D:2:ASN:C	1:D:2:ASN:HD22	2.21	0.43
1:D:75:ASN:HA	1:D:78:LYS:HG3	2.00	0.43
1:A:74:THR:O	1:A:77:GLU:HB2	2.18	0.43
1:A:96:GLN:O	1:A:100:ILE:HG12	2.16	0.43
1:E:2:ASN:O	1:E:5:ASP:N	2.51	0.43
1:A:43:PHE:O	1:A:45:THR:N	2.46	0.43
1:B:168:ILE:HA	1:B:171:ARG:HG3	2.00	0.43
1:D:132:LYS:O	1:D:135:ASN:HB2	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:14:LEU:N	1:B:14:LEU:HD12	2.33	0.43
1:B:144:ASP:OD2	1:B:144:ASP:N	2.50	0.43
1:B:18:ASN:OD1	1:B:18:ASN:N	2.52	0.43
1:B:29:VAL:HG11	1:B:35:SER:HA	2.01	0.43
1:E:163:THR:O	1:E:165:GLU:N	2.48	0.43
1:B:2:ASN:HB3	1:B:5:ASP:HB2	2.00	0.43
1:E:20:TYR:CE1	1:E:46:LYS:HE2	2.55	0.42
1:E:74:THR:HG1	1:E:77:GLU:HG3	1.84	0.42
1:A:115:LYS:H	1:A:115:LYS:CD	2.20	0.42
1:A:183:LEU:HA	1:A:183:LEU:HD12	1.80	0.42
1:E:136:LEU:HD12	1:E:137:ASN:ND2	2.34	0.42
1:D:70:ILE:O	1:D:72:ALA:N	2.52	0.42
1:E:81:LEU:HG	1:E:85:LEU:CD2	2.50	0.42
1:E:167:ASN:OD1	1:E:169:ASN:N	2.53	0.42
1:B:100:ILE:O	1:B:104:THR:HG23	2.18	0.42
1:D:34:SER:OG	1:D:35:SER:N	2.51	0.42
1:D:43:PHE:O	1:D:44:LYS:HB2	2.19	0.42
1:A:47:GLU:OE2	1:A:112:ILE:HD11	2.19	0.42
1:D:36:LYS:O	1:D:40:TYR:HD1	2.03	0.42
1:D:59:SER:C	1:D:61:TRP:N	2.71	0.42
1:D:183:LEU:C	1:D:185:GLY:N	2.73	0.42
1:A:90:GLU:HG2	1:A:160:VAL:CG1	2.49	0.42
1:A:132:LYS:O	1:A:136:LEU:HD13	2.20	0.42
1:E:45:THR:HG1	1:E:47:GLU:CD	2.23	0.42
1:D:59:SER:O	1:D:61:TRP:N	2.52	0.42
1:D:64:GLN:O	1:D:64:GLN:HG2	2.19	0.42
1:D:90:GLU:H	1:D:90:GLU:HG3	1.64	0.42
1:E:168:ILE:O	1:E:172:ILE:HG13	2.20	0.42
1:B:174:LEU:HB3	1:A:151:ILE:HD13	2.01	0.42
1:A:81:LEU:HD12	1:A:81:LEU:O	2.20	0.42
1:B:163:THR:O	1:B:171:ARG:HD2	2.19	0.41
1:E:103:TYR:CD2	1:E:116:MET:HG2	2.55	0.41
1:B:6:LYS:O	1:B:10:VAL:HG23	2.20	0.41
1:A:47:GLU:H	1:A:47:GLU:HG3	1.42	0.41
1:A:76:ARG:HG3	1:A:183:LEU:CD2	2.40	0.41
1:A:167:ASN:OD1	1:A:170:GLU:CB	2.68	0.41
1:B:151:ILE:HD13	1:A:174:LEU:HB3	2.01	0.41
1:D:32:SER:O	1:D:33:GLU:HB2	2.20	0.41
1:E:29:VAL:HG21	1:E:36:LYS:HA	2.02	0.41
1:B:66:LYS:HD3	1:B:69:GLN:NE2	2.35	0.41
1:D:150:LYS:HE3	1:D:150:LYS:HB2	1.82	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:140:TRP:CZ3	1:B:183:LEU:HD22	2.56	0.41
1:E:124:ILE:HD13	1:E:150:LYS:HG2	2.03	0.41
1:D:84:GLU:HG3	1:D:176:ASN:HD21	1.84	0.41
1:E:158:GLY:O	1:E:162:PHE:HB2	2.21	0.41
1:E:87:LEU:HD11	1:E:175:MET:HB2	2.03	0.41
1:E:93:TYR:N	1:E:94:PRO:CD	2.84	0.41
1:E:100:ILE:HG23	1:E:101:GLU:N	2.35	0.41
1:E:131:PHE:HB2	1:E:145:VAL:HG13	2.03	0.41
1:D:40:TYR:O	1:D:44:LYS:N	2.54	0.41
1:E:47:GLU:CD	1:E:47:GLU:H	2.24	0.41
1:B:23:THR:O	1:B:46:LYS:NZ	2.44	0.40
1:B:39:LEU:C	1:B:43:PHE:O	2.60	0.40
1:A:159:ILE:HB	1:A:175:MET:HE1	2.02	0.40
1:E:65:TRP:CD2	1:E:82:TYR:HD1	2.40	0.40
1:E:81:LEU:O	1:E:85:LEU:HD22	2.22	0.40
1:E:181:ILE:O	1:E:184:ASN:HB3	2.20	0.40
1:E:163:THR:C	1:E:165:GLU:N	2.74	0.40
1:B:15:PHE:CD1	1:B:50:PHE:HD2	2.38	0.40
1:D:29:VAL:HG13	1:D:34:SER:C	2.42	0.40
1:A:95:LEU:HD23	1:A:95:LEU:HA	1.86	0.40
1:E:15:PHE:CE1	1:E:50:PHE:HD2	2.39	0.40
1:D:156:VAL:HG13	1:D:175:MET:HE3	2.03	0.40
1:B:12:LYS:HE3	1:B:57:GLU:OE2	2.22	0.40
1:B:122:LYS:HB2	1:B:122:LYS:NZ	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	184/194 (95%)	159 (86%)	18 (10%)	7 (4%)	3 10

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	B	176/194 (91%)	140 (80%)	30 (17%)	6 (3%)	3	13
1	D	180/194 (93%)	145 (81%)	25 (14%)	10 (6%)	2	5
1	E	184/194 (95%)	157 (85%)	23 (12%)	4 (2%)	6	22
All	All	724/776 (93%)	601 (83%)	96 (13%)	27 (4%)	3	11

All (27) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	89	THR
1	A	165	GLU
1	B	17	LYS
1	D	60	LYS
1	D	93	TYR
1	D	143	ASN
1	A	136	LEU
1	A	164	HIS
1	A	166	GLN
1	E	92	TYR
1	E	166	GLN
1	B	67	LYS
1	D	3	LEU
1	D	71	LYS
1	D	92	TYR
1	D	137	ASN
1	A	111	SER
1	E	143	ASN
1	E	164	HIS
1	B	167	ASN
1	D	30	LYS
1	A	104	THR
1	D	44	LYS
1	B	38	ASN
1	B	110	ASN
1	D	29	VAL
1	A	145	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	165/177 (93%)	147 (89%)	18 (11%)	6	19
1	B	158/177 (89%)	143 (90%)	15 (10%)	8	25
1	D	151/177 (85%)	138 (91%)	13 (9%)	10	30
1	E	165/177 (93%)	158 (96%)	7 (4%)	30	63
All	All	639/708 (90%)	586 (92%)	53 (8%)	11	32

All (53) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	5	ASP
1	B	6	LYS
1	B	18	ASN
1	B	58	GLU
1	B	70	ILE
1	B	85	LEU
1	B	88	THR
1	B	107	TYR
1	B	109	THR
1	B	140	TRP
1	B	143	ASN
1	B	157	ASN
1	B	165	GLU
1	B	167	ASN
1	B	180	GLN
1	D	2	ASN
1	D	42	HIS
1	D	45	THR
1	D	47	GLU
1	D	70	ILE
1	D	90	GLU
1	D	96	GLN
1	D	104	THR
1	D	110	ASN
1	D	140	TRP
1	D	166	GLN
1	D	167	ASN
1	D	177	LYS
1	A	47	GLU
1	A	49	LEU

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Mol	Chain	Res	Type
1	A	54	LEU
1	A	64	GLN
1	A	85	LEU
1	A	90	GLU
1	A	105	GLU
1	A	107	TYR
1	A	108	LYS
1	A	115	LYS
1	A	116	MET
1	A	136	LEU
1	A	143	ASN
1	A	167	ASN
1	A	169	ASN
1	A	171	ARG
1	A	173	LYS
1	A	183	LEU
1	E	24	THR
1	E	45	THR
1	E	47	GLU
1	E	54	LEU
1	E	104	THR
1	E	133	GLU
1	E	166	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (52) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	21	ASN
1	B	62	GLN
1	B	64	GLN
1	B	69	GLN
1	B	75	ASN
1	B	83	ASN
1	B	96	GLN
1	B	97	ASN
1	B	113	ASN
1	B	121	ASN
1	B	128	HIS
1	B	154	ASN
1	B	157	ASN
1	B	166	GLN
1	B	167	ASN

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Mol	Chain	Res	Type
1	D	2	ASN
1	D	62	GLN
1	D	75	ASN
1	D	83	ASN
1	D	96	GLN
1	D	110	ASN
1	D	113	ASN
1	D	117	ASN
1	D	143	ASN
1	D	154	ASN
1	D	166	GLN
1	D	167	ASN
1	D	176	ASN
1	A	21	ASN
1	A	38	ASN
1	A	62	GLN
1	A	64	GLN
1	A	69	GLN
1	A	83	ASN
1	A	117	ASN
1	A	121	ASN
1	A	135	ASN
1	A	143	ASN
1	A	157	ASN
1	A	180	GLN
1	A	184	ASN
1	E	18	ASN
1	E	38	ASN
1	E	83	ASN
1	E	96	GLN
1	E	97	ASN
1	E	117	ASN
1	E	121	ASN
1	E	128	HIS
1	E	146	ASN
1	E	157	ASN
1	E	166	GLN

5.3.3 RNA ⓘ

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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

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	D	516	-	4,4,4	0.27	0	6,6,6	0.05	0
2	SO4	A	507	-	4,4,4	0.25	0	6,6,6	0.07	0
2	SO4	A	506	-	4,4,4	0.26	0	6,6,6	0.05	0
2	SO4	E	515	-	4,4,4	0.26	0	6,6,6	0.07	0
2	SO4	E	514	-	4,4,4	0.29	0	6,6,6	0.06	0
2	SO4	A	503	-	4,4,4	0.31	0	6,6,6	0.17	0
2	SO4	E	512	-	4,4,4	0.30	0	6,6,6	0.13	0
2	SO4	A	504	-	4,4,4	0.25	0	6,6,6	0.12	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.

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 ⓘ

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å ²)	Q<0.9
1	A	186/194 (95%)	-0.00	4 (2%)	62 52	28, 47, 98, 130	0
1	B	182/194 (93%)	0.13	5 (2%)	54 44	38, 75, 136, 168	0
1	D	184/194 (94%)	-0.01	3 (1%)	72 66	42, 76, 109, 119	0
1	E	186/194 (95%)	0.15	5 (2%)	54 44	35, 60, 95, 107	0
All	All	738/776 (95%)	0.07	17 (2%)	60 51	28, 65, 113, 168	0

All (17) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	142	ILE	4.0
1	B	110	ASN	3.3
1	B	112	ILE	3.1
1	B	136	LEU	2.6
1	B	107	TYR	2.6
1	A	112	ILE	2.5
1	B	113	ASN	2.4
1	E	145	VAL	2.3
1	E	123	TYR	2.3
1	D	163	THR	2.2
1	D	112	ILE	2.2
1	E	140	TRP	2.1
1	D	140	TRP	2.1
1	A	142	ILE	2.0
1	E	141	SER	2.0
1	A	186	LEU	2.0
1	A	164	HIS	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides 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(\AA^2)	Q<0.9
2	SO4	E	515	5/5	0.70	0.34	165,165,166,167	0
2	SO4	D	516	5/5	0.71	0.22	185,185,187,187	0
2	SO4	E	514	5/5	0.77	0.31	139,142,143,146	0
2	SO4	A	507	5/5	0.77	0.19	142,143,143,143	0
2	SO4	A	506	5/5	0.82	0.38	172,172,173,174	0
2	SO4	A	504	5/5	0.91	0.17	99,102,109,109	0
2	SO4	A	503	5/5	0.94	0.18	96,97,103,104	0
2	SO4	E	512	5/5	0.94	0.13	98,99,103,105	0

6.5 Other polymers [i](#)

There are no such residues in this entry.