



Full wwPDB X-ray Structure Validation Report ⓘ

May 16, 2020 – 01:30 pm BST

PDB ID : 2WSQ
Title : MonoTIM mutant RMM0-1, dimeric form.
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Deposited on : 2009-09-08
Resolution : 2.10 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 : 2.11
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) : 2.11

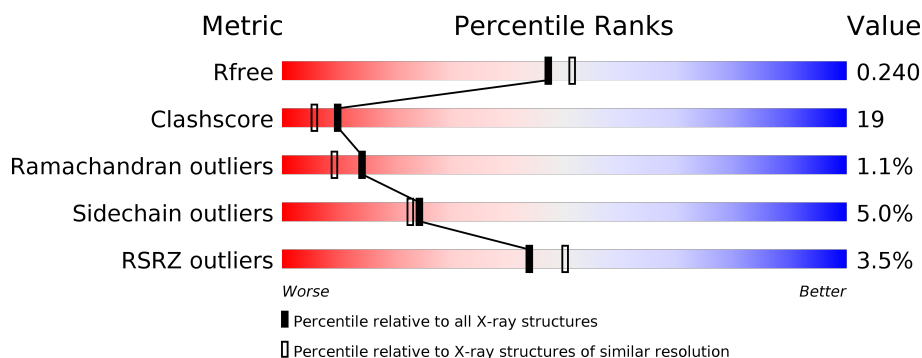
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (2.10-2.10)

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 ≥ 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	242	<div> <div>3%</div> <div> <div></div> <div>71%</div> <div>25%</div> <div>5%</div> </div> </div>
1	B	242	<div> <div>5%</div> <div> <div></div> <div>68%</div> <div>29%</div> <div>.</div> </div> </div>
1	C	242	<div> <div>2%</div> <div> <div></div> <div>70%</div> <div>27%</div> <div>.</div> </div> </div>
1	D	242	<div> <div>3%</div> <div> <div></div> <div>68%</div> <div>28%</div> <div>.</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 8904 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 TRIOSE PHOSPHATE ISOMERASE, GLYCOSOMAL.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	242	Total	C	N	O	S	0	0	0
			1828	1158	324	341	5			
1	B	242	Total	C	N	O	S	0	0	0
			1828	1158	324	341	5			
1	C	242	Total	C	N	O	S	0	0	0
			1828	1158	324	341	5			
1	D	242	Total	C	N	O	S	0	0	0
			1827	1158	324	340	5			

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	43	PRO	SER	engineered mutation	UNP P04789
A	44	SER	THR	engineered mutation	UNP P04789
A	186	THR	ALA	engineered mutation	UNP P04789
B	43	PRO	SER	engineered mutation	UNP P04789
B	44	SER	THR	engineered mutation	UNP P04789
B	186	THR	ALA	engineered mutation	UNP P04789
C	43	PRO	SER	engineered mutation	UNP P04789
C	44	SER	THR	engineered mutation	UNP P04789
C	186	THR	ALA	engineered mutation	UNP P04789
D	43	PRO	SER	engineered mutation	UNP P04789
D	44	SER	THR	engineered mutation	UNP P04789
D	186	THR	ALA	engineered mutation	UNP P04789

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



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

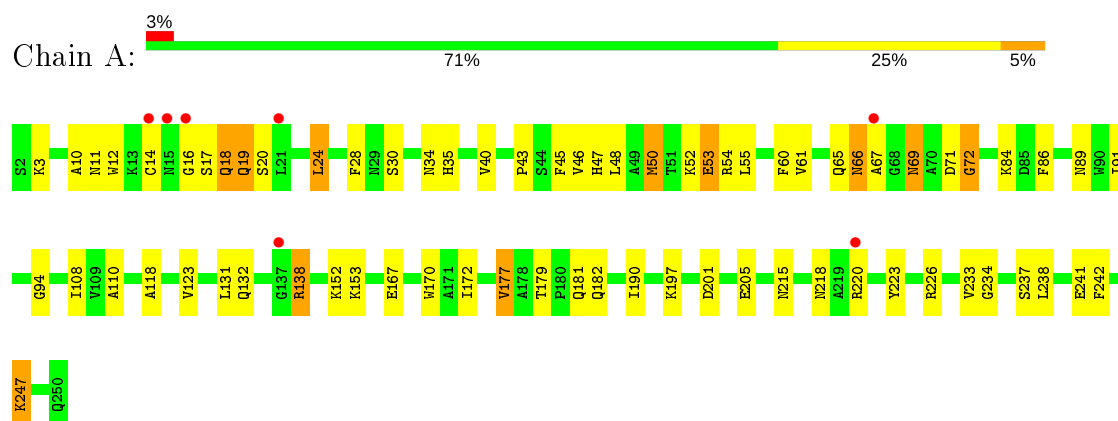
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	376	Total	O	0	0
			376	376		
3	B	370	Total	O	0	0
			370	370		
3	C	373	Total	O	0	0
			373	373		
3	D	434	Total	O	0	0
			434	434		

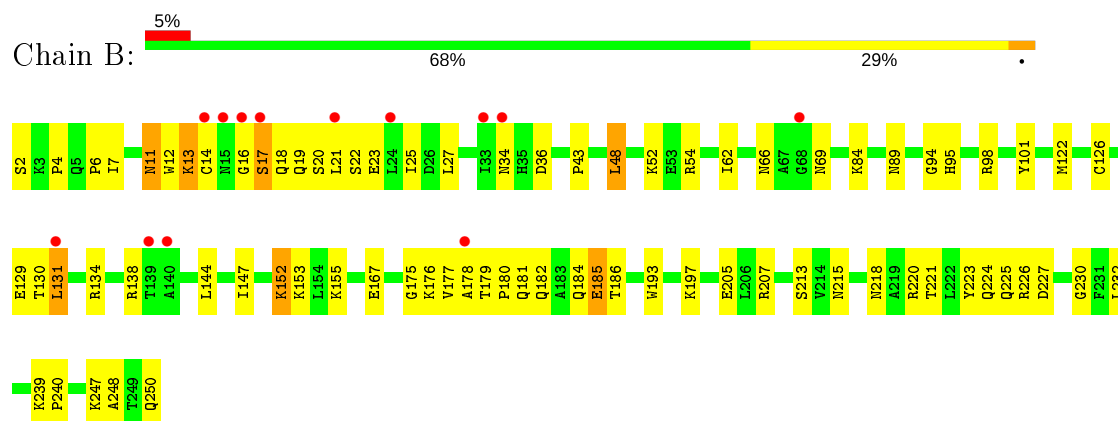
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA 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.

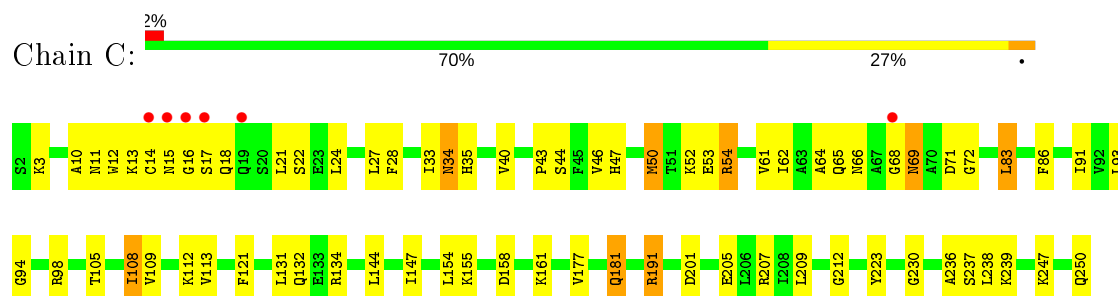
• Molecule 1: TRIOSE PHOSPHATE ISOMERASE, GLYCOSOMAL



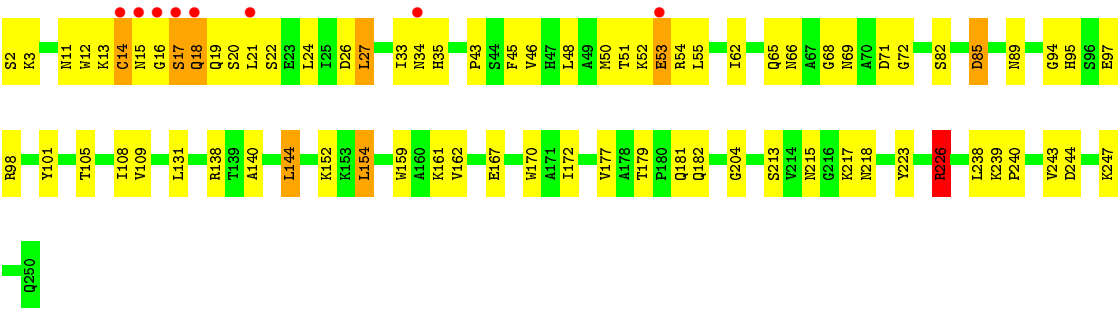
• Molecule 1: TRIOSE PHOSPHATE ISOMERASE, GLYCOSOMAL



• Molecule 1: TRIOSE PHOSPHATE ISOMERASE, GLYCOSOMAL



● Molecule 1: TRIOSE PHOSPHATE ISOMERASE, GLYCOSOMAL



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	90.88Å 94.08Å 108.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	71.25 – 2.10 47.12 – 2.10	Depositor EDS
% Data completeness (in resolution range)	97.8 (71.25-2.10) 97.8 (47.12-2.10)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.01 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.204 , 0.244 0.201 , 0.240	Depositor DCC
R_{free} test set	5446 reflections (10.09%)	wwPDB-VP
Wilson B-factor (Å ²)	31.2	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 62.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	0.008 for k,h,-l	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8904	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 10.19% 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

5.1 Standard geometry

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	0.93	1/1861 (0.1%)	0.79	1/2525 (0.0%)
1	B	0.95	0/1861	0.79	0/2525
1	C	0.93	1/1861 (0.1%)	0.83	3/2525 (0.1%)
1	D	0.99	1/1860 (0.1%)	0.80	1/2524 (0.0%)
All	All	0.95	3/7443 (0.0%)	0.80	5/10099 (0.0%)

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	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	72	GLY	C-N	9.77	1.56	1.34
1	A	72	GLY	C-N	9.15	1.55	1.34
1	C	72	GLY	C-N	8.09	1.52	1.34

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	72	GLY	O-C-N	-7.16	111.24	122.70
1	A	72	GLY	O-C-N	-7.12	111.31	122.70
1	D	226	ARG	NE-CZ-NH1	-5.69	117.45	120.30
1	C	191	ARG	NE-CZ-NH2	-5.55	117.53	120.30
1	C	134	ARG	NE-CZ-NH1	-5.08	117.76	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	72	GLY	Mainchain

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	1828	0	1851	71	0
1	B	1828	0	1851	80	0
1	C	1828	0	1851	62	0
1	D	1827	0	1848	80	0
2	A	5	0	0	0	0
2	B	15	0	0	2	0
2	C	10	0	0	0	0
2	D	10	0	0	0	0
3	A	376	0	0	19	0
3	B	370	0	0	28	0
3	C	373	0	0	14	0
3	D	434	0	0	25	0
All	All	8904	0	7401	283	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 19.

All (283) 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:233:VAL:HG13	1:A:237:SER:HB2	1.30	1.13
1:D:14:CYS:SG	1:D:17:SER:HB3	1.92	1.09
1:A:132:GLN:NE2	1:D:18:GLN:HB2	1.67	1.08
1:A:220:ARG:HG2	3:A:2345:HOH:O	1.58	1.04
1:C:54:ARG:HH11	1:C:54:ARG:HB3	1.22	1.03
1:D:55:LEU:HB2	3:D:2152:HOH:O	1.58	1.01
1:A:190:ILE:HD11	3:A:2238:HOH:O	1.65	0.95
1:A:234:GLY:O	1:A:237:SER:HB3	1.66	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:152:LYS:O	1:B:176:LYS:HE2	1.77	0.84
1:B:52:LYS:HA	1:B:62:ILE:HD11	1.59	0.84
1:D:66:ASN:HD21	1:D:94:GLY:H	1.23	0.84
1:C:54:ARG:HH11	1:C:54:ARG:CB	1.92	0.82
3:A:2240:HOH:O	1:D:46:VAL:HG21	1.77	0.82
1:A:132:GLN:HE22	1:D:18:GLN:HB2	1.42	0.80
1:A:10:ALA:HB1	1:A:237:SER:OG	1.80	0.80
1:D:15:ASN:C	1:D:17:SER:H	1.82	0.80
1:D:17:SER:HB2	1:D:21:LEU:HG	1.64	0.79
1:B:184:GLN:HE22	1:B:227:ASP:HB2	1.47	0.79
1:A:181:GLN:HG2	3:A:2301:HOH:O	1.81	0.78
1:C:17:SER:HB2	1:C:21:LEU:HG	1.66	0.78
1:B:182:GLN:O	3:B:2298:HOH:O	2.02	0.77
1:A:132:GLN:NE2	1:D:18:GLN:CB	2.48	0.77
1:D:154:LEU:HD13	3:D:2283:HOH:O	1.83	0.76
1:D:179:THR:H	1:D:182:GLN:HE21	1.33	0.76
1:D:14:CYS:SG	1:D:17:SER:CB	2.74	0.73
1:B:54:ARG:NH1	1:B:54:ARG:HB3	2.04	0.73
1:B:178:ALA:HA	3:B:2296:HOH:O	1.90	0.72
1:B:66:ASN:HD21	1:B:94:GLY:H	1.37	0.72
1:B:218:ASN:HA	3:B:2336:HOH:O	1.90	0.71
1:A:132:GLN:HE22	1:D:18:GLN:CB	2.03	0.71
1:C:46:VAL:HG22	3:C:2122:HOH:O	1.90	0.70
2:B:1252:SO4:O4	3:B:2368:HOH:O	2.08	0.70
1:A:132:GLN:HB3	3:A:2242:HOH:O	1.92	0.70
1:C:18:GLN:HB2	3:C:2042:HOH:O	1.92	0.70
1:D:52:LYS:HA	1:D:62:ILE:CD1	2.22	0.70
1:D:161:LYS:HE3	3:D:2330:HOH:O	1.91	0.69
1:A:179:THR:H	1:A:182:GLN:HE21	1.36	0.69
1:A:110:ALA:HB1	1:A:153:LYS:HD2	1.75	0.69
1:C:205:GLU:HG3	3:C:2325:HOH:O	1.93	0.68
1:A:30:SER:HA	3:A:2087:HOH:O	1.93	0.67
1:A:24:LEU:HD11	1:A:238:LEU:HA	1.77	0.67
1:C:14:CYS:SG	1:C:17:SER:HB3	2.35	0.67
1:A:53:GLU:HG3	1:A:54:ARG:HG2	1.74	0.67
1:B:101:TYR:HA	1:C:44:SER:HB2	1.78	0.66
1:D:66:ASN:ND2	1:D:94:GLY:H	1.93	0.66
1:B:220:ARG:HB3	3:B:2337:HOH:O	1.96	0.66
1:B:177:VAL:HG22	1:B:213:SER:HB2	1.77	0.66
1:C:105:THR:OG1	1:C:108:ILE:HG22	1.97	0.65
1:B:43:PRO:HG2	1:B:48:LEU:HD12	1.78	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:126:CYS:HB3	3:B:2207:HOH:O	1.96	0.64
1:A:108:ILE:HD13	3:A:2228:HOH:O	1.97	0.64
1:B:184:GLN:HE22	1:B:227:ASP:CB	2.11	0.64
1:B:2:SER:HB3	3:B:2001:HOH:O	1.97	0.64
1:D:82:SER:O	1:D:85:ASP:HB2	1.97	0.64
1:B:131:LEU:HD12	1:C:16:GLY:HA2	1.79	0.64
1:C:236:ALA:HA	1:C:239:LYS:HE2	1.80	0.64
1:D:247:LYS:HE2	3:D:2422:HOH:O	1.98	0.63
1:B:52:LYS:HA	1:B:62:ILE:CD1	2.28	0.63
1:C:54:ARG:HB3	1:C:54:ARG:NH1	2.05	0.63
1:C:12:TRP:HZ2	1:C:28:PHE:HE2	1.47	0.62
1:C:35:HIS:CE1	3:C:2099:HOH:O	2.52	0.62
1:B:177:VAL:HG22	1:B:213:SER:CB	2.29	0.62
1:D:45:PHE:HA	1:D:48:LEU:HD13	1.80	0.62
1:D:179:THR:H	1:D:182:GLN:NE2	1.96	0.62
1:A:220:ARG:NH2	3:A:2344:HOH:O	2.32	0.61
1:A:17:SER:CB	1:A:20:SER:HB3	2.30	0.61
1:D:15:ASN:C	1:D:17:SER:N	2.53	0.61
1:D:26:ASP:HB3	3:D:2082:HOH:O	2.00	0.61
1:A:84:LYS:HD2	1:A:89:ASN:HD22	1.66	0.61
1:A:16:GLY:HA2	1:D:131:LEU:HD22	1.81	0.61
1:B:21:LEU:O	1:B:25:ILE:HG13	2.01	0.61
1:D:22:SER:HB2	3:D:2162:HOH:O	2.00	0.61
1:D:95:HIS:HB2	1:D:109:VAL:HG21	1.81	0.60
1:A:138:ARG:HA	3:A:2253:HOH:O	2.01	0.60
1:A:66:ASN:HD21	1:A:94:GLY:H	1.49	0.60
1:D:53:GLU:HG2	1:D:54:ARG:N	2.16	0.60
1:C:66:ASN:HD21	1:C:94:GLY:H	1.48	0.60
1:C:238:LEU:N	1:C:238:LEU:HD22	2.16	0.60
1:D:179:THR:HG22	3:D:2348:HOH:O	2.01	0.60
1:A:45:PHE:HA	1:A:48:LEU:HD13	1.85	0.59
1:C:68:GLY:O	1:C:69:ASN:HB2	2.01	0.59
1:D:18:GLN:H	1:D:21:LEU:HG	1.68	0.59
1:D:204:GLY:HA3	3:D:2371:HOH:O	2.01	0.58
1:B:226:ARG:NH2	3:B:2344:HOH:O	2.34	0.58
1:B:184:GLN:NE2	1:B:227:ASP:HB2	2.17	0.58
1:A:131:LEU:HD13	1:A:131:LEU:C	2.24	0.58
1:D:51:THR:HG23	3:D:2152:HOH:O	2.04	0.58
1:C:52:LYS:HD2	1:C:86:PHE:CE2	2.39	0.58
1:D:53:GLU:HG2	1:D:54:ARG:HG2	1.85	0.57
1:C:12:TRP:O	1:C:13:LYS:HG3	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:18:GLN:O	1:C:22:SER:HB2	2.05	0.57
1:A:18:GLN:NE2	1:A:50:MET:HG3	2.20	0.57
1:A:167:GLU:HG2	1:A:172:ILE:HD11	1.87	0.57
1:C:112:LYS:HE2	3:C:2164:HOH:O	2.05	0.56
1:A:17:SER:HB2	1:A:20:SER:HB3	1.87	0.56
1:D:52:LYS:HA	1:D:62:ILE:HD11	1.86	0.56
1:B:12:TRP:HB2	1:B:43:PRO:HA	1.88	0.56
1:C:34:ASN:HD21	1:C:250:GLN:HE22	1.53	0.56
1:D:97:GLU:HB3	1:D:101:TYR:CE2	2.41	0.55
1:A:132:GLN:HE21	1:D:18:GLN:HB2	1.67	0.55
1:D:177:VAL:HG21	1:D:218:ASN:ND2	2.21	0.55
1:B:197:LYS:HG2	3:B:2309:HOH:O	2.06	0.55
1:C:11:ASN:OD1	1:C:65:GLN:OE1	2.25	0.55
1:A:52:LYS:NZ	3:A:2129:HOH:O	2.13	0.55
1:C:131:LEU:HD12	1:C:132:GLN:NE2	2.22	0.55
1:D:27:LEU:O	1:D:27:LEU:HD23	2.07	0.55
1:B:54:ARG:HH11	1:B:54:ARG:HB3	1.70	0.55
1:D:46:VAL:HG22	1:D:46:VAL:O	2.07	0.54
1:B:16:GLY:N	3:B:2053:HOH:O	2.40	0.54
1:B:19:GLN:HG3	3:B:2074:HOH:O	2.08	0.54
1:D:13:LYS:HE3	3:D:2410:HOH:O	2.07	0.53
1:A:153:LYS:HE2	3:A:2225:HOH:O	2.08	0.53
1:B:184:GLN:NE2	1:B:227:ASP:OD2	2.41	0.53
1:A:110:ALA:CB	1:A:153:LYS:HD2	2.38	0.53
1:C:161:LYS:HE2	3:C:2288:HOH:O	2.08	0.53
1:D:24:LEU:HD11	1:D:238:LEU:HA	1.90	0.53
1:B:11:ASN:HD22	1:B:232:LEU:HD11	1.74	0.53
1:A:18:GLN:O	1:A:19:GLN:HB2	2.09	0.52
1:A:14:CYS:HA	3:A:2043:HOH:O	2.09	0.52
1:D:105:THR:HG23	3:D:2276:HOH:O	2.09	0.52
1:A:226:ARG:NH1	3:A:2351:HOH:O	2.42	0.52
1:D:15:ASN:O	1:D:17:SER:N	2.43	0.52
1:B:84:LYS:HZ2	1:B:89:ASN:HD22	1.58	0.51
1:A:69:ASN:OD1	1:A:71:ASP:HB2	2.10	0.51
1:D:98:ARG:NE	3:D:2263:HOH:O	2.43	0.51
1:B:52:LYS:HG3	1:B:62:ILE:HD12	1.92	0.51
1:D:159:TRP:HA	1:D:162:VAL:HG22	1.93	0.51
1:B:220:ARG:HB2	3:B:2327:HOH:O	2.10	0.51
1:B:122:MET:HG2	3:B:2254:HOH:O	2.10	0.51
1:D:12:TRP:HB2	1:D:43:PRO:HA	1.93	0.51
1:A:11:ASN:ND2	1:A:65:GLN:OE1	2.45	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:68:GLY:HA3	3:D:2262:HOH:O	2.11	0.50
1:D:52:LYS:CA	1:D:62:ILE:CD1	2.90	0.50
1:B:144:LEU:O	1:B:147:ILE:HG22	2.10	0.50
1:A:55:LEU:HD21	1:A:60:PHE:HB3	1.93	0.50
1:A:67:ALA:HB1	3:D:2271:HOH:O	2.10	0.50
1:A:84:LYS:HD2	1:A:89:ASN:ND2	2.26	0.50
1:A:67:ALA:CB	3:D:2271:HOH:O	2.59	0.50
1:B:129:GLU:OE2	1:B:134:ARG:HB2	2.10	0.50
1:B:134:ARG:HA	3:B:2266:HOH:O	2.12	0.50
1:B:11:ASN:ND2	1:B:232:LEU:HD11	2.27	0.50
1:D:244:ASP:CG	3:D:2418:HOH:O	2.50	0.50
1:C:91:ILE:HG23	1:C:121:PHE:CD2	2.47	0.50
1:B:13:LYS:HB3	3:B:2039:HOH:O	2.12	0.49
1:B:16:GLY:O	1:B:17:SER:HB3	2.12	0.49
1:B:6:PRO:HD2	1:B:36:ASP:O	2.13	0.49
1:B:19:GLN:O	1:B:23:GLU:HG3	2.13	0.49
1:A:215:ASN:HB2	1:A:241:GLU:OE2	2.13	0.49
1:B:179:THR:HG23	1:B:182:GLN:NE2	2.28	0.49
1:C:34:ASN:HD21	1:C:250:GLN:NE2	2.10	0.49
1:C:132:GLN:HB3	3:C:2256:HOH:O	2.12	0.49
1:B:138:ARG:NH1	3:B:2265:HOH:O	2.46	0.49
1:C:181:GLN:H	1:C:181:GLN:CD	2.16	0.48
1:D:226:ARG:HG3	1:D:226:ARG:NH1	2.28	0.48
1:D:33:ILE:HA	1:D:247:LYS:HZ2	1.78	0.48
1:D:108:ILE:HD12	3:D:2276:HOH:O	2.12	0.48
1:C:40:VAL:HG22	1:C:61:VAL:HG23	1.96	0.48
1:D:177:VAL:HG23	3:D:2341:HOH:O	2.11	0.48
1:A:131:LEU:HB2	1:A:170:TRP:HB3	1.96	0.48
1:D:179:THR:N	1:D:182:GLN:HE21	2.08	0.48
1:A:12:TRP:HZ2	1:A:28:PHE:HE2	1.60	0.47
1:A:50:MET:HB3	3:A:2060:HOH:O	2.13	0.47
1:D:138:ARG:NH2	3:D:2308:HOH:O	2.47	0.47
1:B:27:LEU:HD11	1:B:240:PRO:HB3	1.95	0.47
1:C:105:THR:O	1:C:108:ILE:HG23	2.14	0.47
1:C:27:LEU:HD11	3:C:2085:HOH:O	2.13	0.47
1:C:50:MET:O	1:C:53:GLU:HG2	2.13	0.47
1:D:177:VAL:HG22	1:D:213:SER:O	2.14	0.47
1:A:12:TRP:HB2	1:A:43:PRO:HA	1.95	0.47
1:A:40:VAL:HG22	1:A:61:VAL:CG2	2.44	0.47
1:C:12:TRP:HZ2	1:C:28:PHE:CE2	2.31	0.47
1:A:24:LEU:HD11	1:A:238:LEU:CA	2.42	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:SER:O	1:B:54:ARG:NH2	2.48	0.47
1:B:247:LYS:HG3	3:B:2361:HOH:O	2.14	0.47
1:C:191:ARG:HD2	3:C:2355:HOH:O	2.15	0.47
1:D:52:LYS:HA	1:D:62:ILE:HD12	1.98	0.46
1:A:197:LYS:HG2	3:A:2312:HOH:O	2.16	0.46
1:B:84:LYS:NZ	1:B:89:ASN:HD22	2.13	0.46
1:B:220:ARG:O	1:B:224:GLN:HG3	2.15	0.46
1:D:239:LYS:HB3	1:D:240:PRO:HD2	1.98	0.46
1:D:243:VAL:O	1:D:247:LYS:HG2	2.16	0.46
1:A:34:ASN:OD1	1:A:35:HIS:N	2.38	0.46
1:A:43:PRO:HG2	1:A:48:LEU:HD12	1.98	0.46
1:B:247:LYS:O	1:B:250:GLN:HG2	2.14	0.46
1:A:233:VAL:HG11	1:A:242:PHE:CE1	2.51	0.46
1:B:185:GLU:CG	1:B:186:THR:N	2.79	0.46
1:C:14:CYS:SG	1:C:17:SER:CB	3.04	0.46
1:A:118:ALA:HB3	3:A:2235:HOH:O	2.16	0.46
1:C:3:LYS:NZ	1:C:223:TYR:O	2.48	0.46
1:B:19:GLN:HG2	1:B:20:SER:H	1.81	0.46
1:D:2:SER:HB2	3:D:2003:HOH:O	2.16	0.46
1:C:105:THR:O	1:C:108:ILE:CG2	2.64	0.45
1:B:144:LEU:HA	1:B:147:ILE:HG22	1.98	0.45
1:D:154:LEU:HD21	1:D:162:VAL:HG11	1.97	0.45
1:C:12:TRP:HB2	1:C:43:PRO:HA	1.98	0.45
1:D:131:LEU:HB2	1:D:170:TRP:HB3	1.98	0.45
1:A:40:VAL:HG22	1:A:61:VAL:HG23	1.99	0.45
1:A:17:SER:HA	3:A:2055:HOH:O	2.16	0.45
1:A:247:LYS:N	1:A:247:LYS:HD2	2.32	0.45
1:B:98:ARG:NH1	3:B:2215:HOH:O	2.49	0.45
1:C:91:ILE:HD12	1:C:93:LEU:HD21	1.99	0.45
1:B:193:TRP:CZ2	1:B:197:LYS:HG3	2.52	0.45
1:A:45:PHE:CA	1:A:48:LEU:HD13	2.46	0.44
1:B:129:GLU:HG2	1:B:130:THR:O	2.16	0.44
1:B:6:PRO:HG3	1:B:223:TYR:OH	2.18	0.44
1:B:239:LYS:HE3	3:B:2351:HOH:O	2.16	0.44
1:D:167:GLU:HG2	1:D:172:ILE:HD11	2.00	0.44
1:B:167:GLU:HB3	3:B:2207:HOH:O	2.16	0.44
1:C:12:TRP:CD1	1:C:238:LEU:HD13	2.52	0.44
1:C:209:LEU:HD22	1:C:230:GLY:HA3	1.99	0.44
1:C:53:GLU:HG3	1:C:54:ARG:HG2	2.00	0.44
1:D:66:ASN:HD21	1:D:94:GLY:N	2.03	0.44
1:B:95:HIS:HB3	1:B:126:CYS:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:17:SER:N	3:B:2053:HOH:O	2.45	0.44
1:D:179:THR:HA	3:D:2348:HOH:O	2.16	0.44
1:D:11:ASN:HD21	1:D:65:GLN:NE2	2.16	0.44
1:C:109:VAL:O	1:C:113:VAL:HG23	2.18	0.44
1:B:19:GLN:HG2	1:B:20:SER:N	2.32	0.43
1:B:34:ASN:HB2	3:B:2107:HOH:O	2.16	0.43
1:D:215:ASN:OD1	1:D:217:LYS:HB2	2.18	0.43
1:D:45:PHE:HA	1:D:48:LEU:CD1	2.48	0.43
2:B:1253:SO4:O2	1:C:98:ARG:NH2	2.47	0.43
1:C:10:ALA:HB1	1:C:237:SER:HB2	2.01	0.43
1:C:40:VAL:HG22	1:C:61:VAL:CG2	2.49	0.43
1:A:45:PHE:HA	1:A:48:LEU:CD1	2.48	0.43
1:A:46:VAL:HG21	3:D:2294:HOH:O	2.17	0.43
1:C:54:ARG:CG	1:C:54:ARG:HH11	2.30	0.43
1:D:17:SER:HA	1:D:20:SER:OG	2.19	0.43
1:B:4:PRO:HB2	1:B:207:ARG:HD2	2.01	0.43
1:B:94:GLY:O	1:B:98:ARG:HG3	2.19	0.43
1:D:226:ARG:NE	3:D:2406:HOH:O	2.51	0.43
1:C:47:HIS:HD2	3:C:2120:HOH:O	2.01	0.43
1:D:181:GLN:H	1:D:181:GLN:CD	2.22	0.43
1:A:152:LYS:HE2	1:B:175:GLY:O	2.19	0.43
1:C:154:LEU:HD22	1:C:158:ASP:HB3	1.99	0.43
1:D:140:ALA:O	1:D:144:LEU:HB2	2.19	0.43
1:B:18:GLN:HB3	3:B:2059:HOH:O	2.19	0.42
1:B:180:PRO:O	1:B:225:GLN:HG3	2.20	0.42
1:D:177:VAL:HG22	1:D:213:SER:CB	2.49	0.42
1:B:152:LYS:HB3	1:B:152:LYS:HE3	1.68	0.42
1:D:152:LYS:NZ	3:D:2318:HOH:O	2.41	0.42
1:C:144:LEU:HA	1:C:147:ILE:HG22	2.02	0.42
1:D:50:MET:HA	1:D:53:GLU:OE2	2.19	0.42
1:D:69:ASN:HD22	1:D:71:ASP:HB2	1.83	0.42
1:A:16:GLY:HA2	1:D:131:LEU:CD2	2.49	0.42
1:A:177:VAL:HG21	1:A:218:ASN:ND2	2.35	0.42
1:B:218:ASN:HB2	3:B:2338:HOH:O	2.18	0.42
3:A:2243:HOH:O	1:D:18:GLN:HB3	2.20	0.42
1:B:220:ARG:HA	1:B:248:ALA:HB1	2.01	0.42
1:C:191:ARG:HB3	3:C:2355:HOH:O	2.20	0.42
1:B:23:GLU:HB3	3:B:2072:HOH:O	2.20	0.41
1:C:18:GLN:OE1	1:C:50:MET:HG3	2.20	0.41
1:C:66:ASN:ND2	1:C:94:GLY:H	2.17	0.41
1:A:24:LEU:HD11	1:A:238:LEU:C	2.41	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:7:ILE:O	1:B:230:GLY:HA3	2.21	0.41
1:C:52:LYS:HA	1:C:62:ILE:HD11	2.02	0.41
1:D:247:LYS:CE	3:D:2421:HOH:O	2.67	0.41
1:C:112:LYS:NZ	3:C:2244:HOH:O	2.49	0.41
1:D:226:ARG:HH11	1:D:226:ARG:HG3	1.86	0.41
1:B:19:GLN:CG	1:B:20:SER:N	2.83	0.41
1:C:247:LYS:CE	3:C:2088:HOH:O	2.68	0.41
1:C:64:ALA:CB	1:C:83:LEU:HD21	2.51	0.41
1:D:3:LYS:NZ	1:D:223:TYR:O	2.53	0.41
1:D:34:ASN:OD1	1:D:35:HIS:N	2.53	0.41
1:B:184:GLN:HE22	1:B:227:ASP:CG	2.24	0.41
1:B:193:TRP:CE2	1:B:197:LYS:HG3	2.55	0.41
1:B:184:GLN:OE1	1:B:225:GLN:HB3	2.21	0.41
1:C:24:LEU:O	1:C:27:LEU:HB3	2.19	0.41
1:A:66:ASN:ND2	1:A:94:GLY:H	2.16	0.41
1:C:33:ILE:HG23	3:C:2094:HOH:O	2.20	0.41
1:C:177:VAL:HG23	1:C:212:GLY:O	2.21	0.41
1:A:52:LYS:HD3	1:A:86:PHE:CE2	2.56	0.41
1:B:179:THR:HG23	1:B:182:GLN:HE21	1.84	0.41
1:B:221:THR:HG23	3:B:2337:HOH:O	2.21	0.41
1:A:18:GLN:HG2	1:A:47:HIS:CD2	2.56	0.40
1:A:233:VAL:HG11	1:A:242:PHE:CD1	2.56	0.40
1:A:35:HIS:CE1	3:A:2103:HOH:O	2.74	0.40
1:C:94:GLY:O	1:C:98:ARG:HG3	2.20	0.40
1:B:184:GLN:OE1	1:B:226:ARG:N	2.55	0.40
1:B:215:ASN:ND2	3:B:2326:HOH:O	2.54	0.40
1:A:3:LYS:NZ	1:A:223:TYR:O	2.54	0.40
1:B:138:ARG:N	3:B:2266:HOH:O	2.55	0.40
1:A:91:ILE:HG22	1:A:123:VAL:HG22	2.04	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	240/242 (99%)	226 (94%)	11 (5%)	3 (1%)	12	7
1	B	240/242 (99%)	225 (94%)	13 (5%)	2 (1%)	19	15
1	C	240/242 (99%)	228 (95%)	9 (4%)	3 (1%)	12	7
1	D	240/242 (99%)	228 (95%)	9 (4%)	3 (1%)	12	7
All	All	960/968 (99%)	907 (94%)	42 (4%)	11 (1%)	14	9

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	19	GLN
1	B	14	CYS
1	A	69	ASN
1	B	17	SER
1	D	14	CYS
1	D	16	GLY
1	C	15	ASN
1	A	66	ASN
1	C	34	ASN
1	D	17	SER
1	C	69	ASN

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	190/190 (100%)	181 (95%)	9 (5%)	26	25
1	B	190/190 (100%)	179 (94%)	11 (6%)	20	17
1	C	190/190 (100%)	181 (95%)	9 (5%)	26	25
1	D	189/190 (100%)	180 (95%)	9 (5%)	25	24
All	All	759/760 (100%)	721 (95%)	38 (5%)	24	23

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	18	GLN
1	A	24	LEU
1	A	50	MET
1	A	53	GLU
1	A	138	ARG
1	A	177	VAL
1	A	201	ASP
1	A	205	GLU
1	A	247	LYS
1	B	11	ASN
1	B	13	LYS
1	B	48	LEU
1	B	69	ASN
1	B	131	LEU
1	B	152	LYS
1	B	153	LYS
1	B	155	LYS
1	B	181	GLN
1	B	185	GLU
1	B	205	GLU
1	C	50	MET
1	C	54	ARG
1	C	71	ASP
1	C	83	LEU
1	C	108	ILE
1	C	155	LYS
1	C	181	GLN
1	C	201	ASP
1	C	207	ARG
1	D	18	GLN
1	D	19	GLN
1	D	27	LEU
1	D	53	GLU
1	D	85	ASP
1	D	89	ASN
1	D	144	LEU
1	D	154	LEU
1	D	226	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN

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Mol	Chain	Res	Type
1	A	18	GLN
1	A	66	ASN
1	A	89	ASN
1	A	132	GLN
1	A	181	GLN
1	A	182	GLN
1	A	250	GLN
1	B	65	GLN
1	B	66	ASN
1	B	69	ASN
1	B	89	ASN
1	B	106	ASN
1	B	182	GLN
1	C	34	ASN
1	C	65	GLN
1	C	66	ASN
1	C	69	ASN
1	C	89	ASN
1	C	106	ASN
1	C	132	GLN
1	C	250	GLN
1	D	65	GLN
1	D	66	ASN
1	D	69	ASN
1	D	182	GLN

5.3.3 RNA ⓘ

There are no RNA molecules 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	B	1253	-	4,4,4	0.31	0	6,6,6	0.49	0
2	SO4	C	1251	-	4,4,4	0.22	0	6,6,6	0.41	0
2	SO4	C	1252	-	4,4,4	0.20	0	6,6,6	0.42	0
2	SO4	A	1251	-	4,4,4	0.33	0	6,6,6	0.29	0
2	SO4	D	1252	-	4,4,4	0.23	0	6,6,6	0.60	0
2	SO4	B	1251	-	4,4,4	0.28	0	6,6,6	0.45	0
2	SO4	B	1252	-	4,4,4	0.43	0	6,6,6	0.30	0
2	SO4	D	1251	-	4,4,4	0.39	0	6,6,6	0.29	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.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1253	SO4	1	0
2	B	1252	SO4	1	0

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, 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	242/242 (100%)	0.05	7 (2%) 51 57	19, 31, 48, 66	0
1	B	242/242 (100%)	0.15	13 (5%) 25 31	18, 33, 52, 75	0
1	C	242/242 (100%)	-0.04	6 (2%) 57 62	16, 30, 48, 71	0
1	D	242/242 (100%)	-0.10	8 (3%) 46 53	16, 26, 47, 68	0
All	All	968/968 (100%)	0.01	34 (3%) 44 50	16, 30, 51, 75	0

All (34) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	17	SER	7.0
1	B	16	GLY	5.9
1	B	15	ASN	5.6
1	D	15	ASN	5.2
1	C	15	ASN	4.7
1	C	14	CYS	4.2
1	A	16	GLY	4.1
1	C	16	GLY	4.1
1	A	15	ASN	4.1
1	D	17	SER	4.0
1	D	16	GLY	3.8
1	D	18	GLN	3.7
1	C	19	GLN	3.4
1	B	68	GLY	3.4
1	B	14	CYS	3.3
1	B	17	SER	3.2
1	B	139	THR	3.2
1	A	137	GLY	3.1
1	B	140	ALA	3.0
1	B	21	LEU	2.8
1	A	67	ALA	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	131	LEU	2.7
1	D	53	GLU	2.6
1	D	14	CYS	2.6
1	A	21	LEU	2.6
1	D	34	ASN	2.3
1	B	178	ALA	2.3
1	A	14	CYS	2.2
1	B	24	LEU	2.2
1	C	68	GLY	2.2
1	B	33	ILE	2.1
1	B	34	ASN	2.1
1	A	220	ARG	2.0
1	D	21	LEU	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. 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	B	1252	5/5	0.81	0.34	68,68,69,70	0
2	SO4	B	1253	5/5	0.92	0.17	60,62,64,65	0
2	SO4	C	1252	5/5	0.97	0.12	55,56,58,58	0
2	SO4	B	1251	5/5	0.98	0.09	34,36,39,40	0
2	SO4	D	1252	5/5	0.99	0.08	20,21,25,25	0
2	SO4	C	1251	5/5	0.99	0.09	31,32,33,34	0
2	SO4	A	1251	5/5	0.99	0.09	34,35,36,38	0
2	SO4	D	1251	5/5	0.99	0.05	29,30,31,33	0

6.5 Other polymers

There are no such residues in this entry.