



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:12 PM GMT

PDB ID : 1NXU
Title : CRYSTAL STRUCTURE OF E. COLI HYPOTHETICAL OXIDOREDUCTASE YIAK NORTHEAST STRUCTURAL GENOMICS CONSORTIUM TARGET ER82.
Authors : Forouhar, F.; Lee, I.; Benach, J.; Kulkarni, K.; Xiao, R.; Acton, T.B.; Shastry, R.; Rost, B.; Montelione, G.T.; Tong, L.; Northeast Structural Genomics Consortium (NESG)
Deposited on : 2003-02-11
Resolution : 1.80 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
<http://wwpdb.org/validation/2016/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 : 1.7 (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

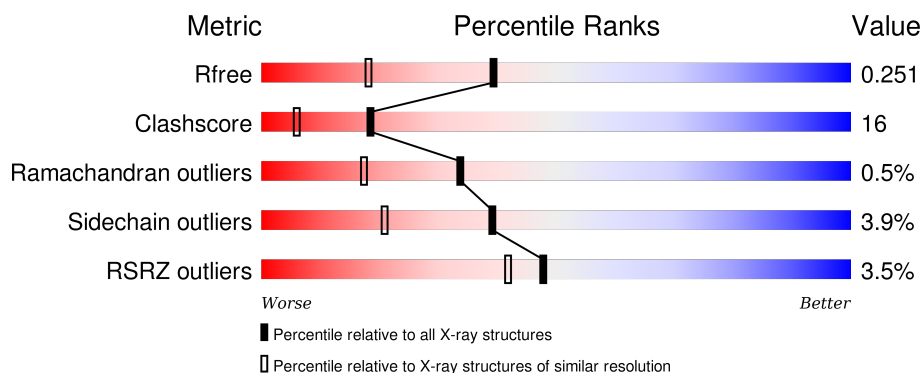
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 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	4533 (1.80-1.80)
Clashscore	102246	5383 (1.80-1.80)
Ramachandran outliers	100387	5320 (1.80-1.80)
Sidechain outliers	100360	5319 (1.80-1.80)
RSRZ outliers	91569	4547 (1.80-1.80)

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 ≥ 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	333	<div> <div>2%</div> <div>74%</div> <div>23%</div> <div>.</div> </div>
1	B	333	<div> <div>4%</div> <div>68%</div> <div>30%</div> <div>.</div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 5813 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 Hypothetical oxidoreductase yiaK.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	332	Total	C	N	O	S	Se	0	0	0
			2562	1604	451	489	3	15			
1	B	333	Total	C	N	O	S	Se	0	0	0
			2567	1607	452	490	3	15			

There are 32 discrepancies between the modelled and reference sequences:

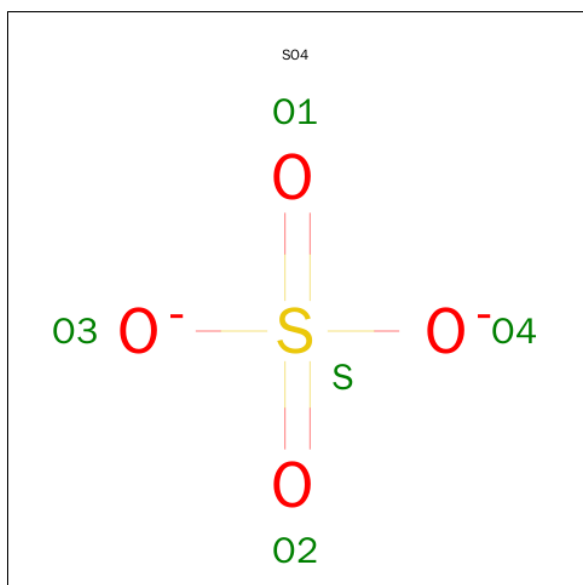
Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	32	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	93	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	94	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	118	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	144	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	170	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	173	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	175	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	177	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	182	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	221	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	229	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	235	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	285	MSE	MET	MODIFIED RESIDUE	UNP P37672
A	333	ALA	-	CLONING ARTIFACT	UNP P37672
B	1	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	32	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	93	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	94	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	118	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	144	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	170	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	173	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	175	MSE	MET	MODIFIED RESIDUE	UNP P37672

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Chain	Residue	Modelled	Actual	Comment	Reference
B	177	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	182	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	221	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	229	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	235	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	285	MSE	MET	MODIFIED RESIDUE	UNP P37672
B	333	ALA	-	CLONING ARTIFACT	UNP P37672

- 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	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		

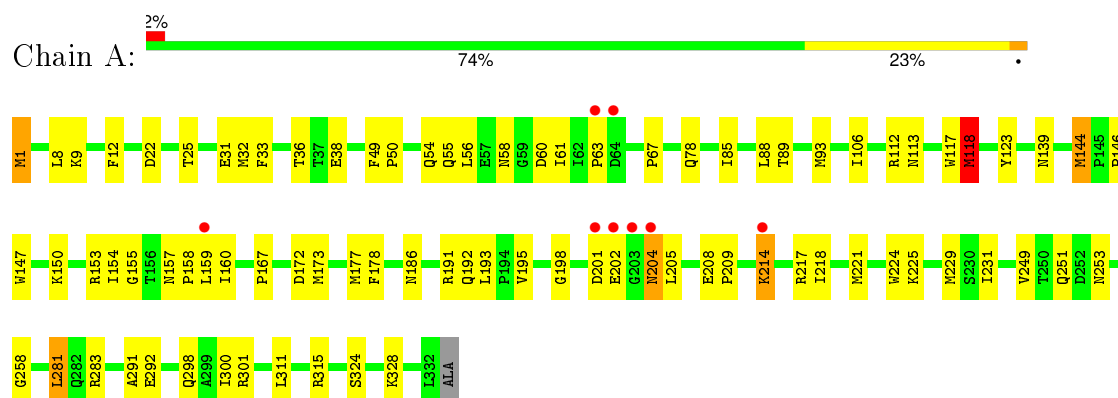
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	335	Total	O	0	0
			335	335		
3	B	329	Total	O	0	0
			329	329		

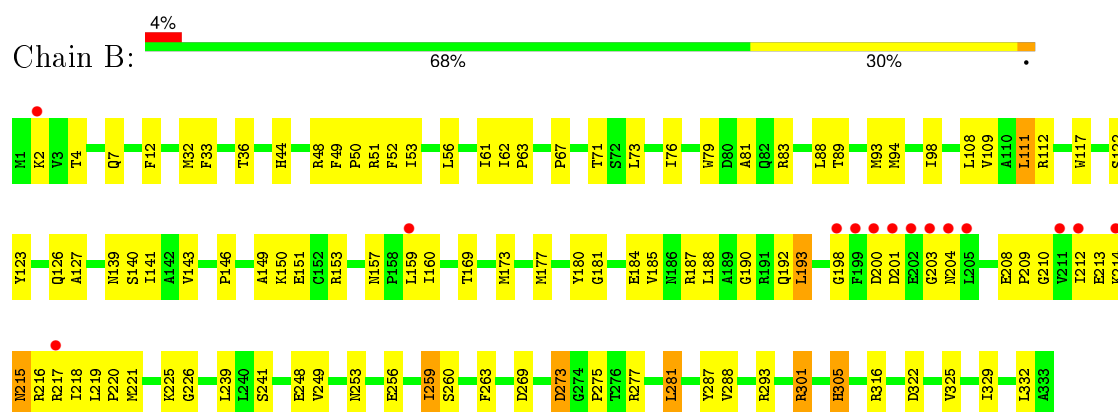
3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Hypothetical oxidoreductase yiaK



• Molecule 1: Hypothetical oxidoreductase yiaK



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	58.01Å 51.23Å 108.93Å 90.00° 103.78° 90.00°	Depositor
Resolution (Å)	37.90 – 1.80 37.90 – 1.79	Depositor EDS
% Data completeness (in resolution range)	68.1 (37.90-1.80) 91.0 (37.90-1.79)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.31 (at 1.79Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.192 , 0.235 0.212 , 0.251	Depositor DCC
R_{free} test set	4191 reflections (7.83%)	DCC
Wilson B-factor (Å ²)	18.7	Xtriage
Anisotropy	0.295	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.26 , 46.2	EDS
Estimated twinning fraction	0.020 for h,-k,-h-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 106854 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5813	wwPDB-VP
Average B, all atoms (Å ²)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.93% 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, 0.375 respectively for untwinned datasets, and 0.333, 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	0.31	0/2594	0.55	0/3487
1	B	0.31	0/2599	0.55	0/3494
All	All	0.31	0/5193	0.55	0/6981

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

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	123	TYR	Sidechain

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	2562	0	2545	82	0
1	B	2567	0	2550	94	0
2	A	10	0	0	0	0
2	B	10	0	0	0	0
3	A	335	0	0	11	0
3	B	329	0	0	5	0
All	All	5813	0	5095	159	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 16.

All (159) 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:B:67:PRO:HB3	1:B:93:MSE:HE2	1.48	0.95
1:B:94:MSE:HE1	1:B:127:ALA:HB2	1.56	0.88
1:B:94:MSE:CE	1:B:127:ALA:HB2	2.08	0.84
1:B:193:LEU:HD13	1:B:209:PRO:HG3	1.64	0.78
1:A:32:MSE:O	1:A:36:THR:HG23	1.86	0.76
1:A:144:MSE:HG2	1:A:218:ILE:HG23	1.69	0.75
1:B:157:ASN:HD22	1:B:157:ASN:H	1.36	0.72
1:A:204:ASN:ND2	1:A:205:LEU:H	1.88	0.72
1:B:109:VAL:CG2	1:B:263:PHE:HB2	2.20	0.71
1:B:67:PRO:HB3	1:B:93:MSE:CE	2.21	0.70
1:A:178:PHE:O	1:A:218:ILE:HD11	1.90	0.70
1:A:154:ILE:HD11	1:B:288:VAL:HG22	1.73	0.69
1:B:198:GLY:HA2	1:B:219:LEU:HB2	1.74	0.69
1:B:32:MSE:O	1:B:36:THR:HG23	1.93	0.69
1:A:291:ALA:HB2	1:B:153:ARG:HG3	1.78	0.65
1:A:225:LYS:HD2	1:B:221:MSE:HG3	1.78	0.65
1:A:229:MSE:HE2	1:B:173:MSE:HE3	1.79	0.64
1:A:281:LEU:HD13	1:B:239:LEU:HD21	1.80	0.63
1:A:192:GLN:HE22	1:A:208:GLU:HG2	1.64	0.62
1:A:144:MSE:HG2	1:A:218:ILE:CG2	2.30	0.61
1:A:9:LYS:HE3	1:A:31:GLU:HG2	1.83	0.61
1:B:109:VAL:HG22	1:B:263:PHE:HB2	1.82	0.61
1:A:36:THR:HG22	3:A:1118:HOH:O	2.01	0.60
1:A:139:ASN:HA	1:A:157:ASN:ND2	2.16	0.60
1:B:301:ARG:HD2	3:B:1293:HOH:O	2.02	0.60
1:B:36:THR:HG22	3:B:1025:HOH:O	2.02	0.60
1:B:200:ASP:OD1	1:B:203:GLY:HA3	2.03	0.59
1:B:140:SER:HB2	1:B:256:GLU:OE2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:157:ASN:H	1:B:157:ASN:ND2	2.00	0.59
1:A:157:ASN:HD22	1:A:157:ASN:H	1.49	0.59
1:B:273:ASP:OD2	1:B:275:PRO:HD2	2.02	0.59
1:A:249:VAL:HG13	1:A:253:ASN:HD22	1.67	0.59
1:A:301:ARG:HH11	1:A:301:ARG:HG2	1.68	0.59
1:B:249:VAL:O	1:B:253:ASN:HB2	2.03	0.58
1:A:251:GLN:HG2	3:B:1271:HOH:O	2.02	0.58
1:B:79:TRP:CE3	1:B:93:MSE:HE1	2.39	0.58
1:B:79:TRP:HB3	1:B:93:MSE:HE1	1.86	0.58
1:A:158:PRO:HG3	3:A:1164:HOH:O	2.03	0.58
1:B:4:THR:HG23	1:B:7:GLN:H	1.69	0.58
1:B:81:ALA:HB2	1:B:111:LEU:HD21	1.86	0.58
1:A:78:GLN:NE2	1:A:112:ARG:HH12	2.02	0.57
1:B:157:ASN:HD22	1:B:157:ASN:N	2.01	0.57
1:A:55:GLN:HG2	3:A:1224:HOH:O	2.03	0.57
1:B:4:THR:HG22	1:B:7:GLN:HB2	1.86	0.56
1:A:56:LEU:HD11	1:A:63:PRO:HG3	1.88	0.56
1:A:49:PHE:HB3	1:A:50:PRO:HD3	1.88	0.55
1:B:215:ASN:HD22	1:B:215:ASN:N	2.05	0.55
1:A:202:GLU:HA	1:A:217:ARG:NH1	2.22	0.55
1:B:49:PHE:HB3	1:B:50:PRO:HD3	1.89	0.54
1:B:184:GLU:HG3	1:B:187:ARG:NH2	2.22	0.54
1:A:177:MSE:HE2	1:A:195:VAL:HG21	1.90	0.54
1:A:292:GLU:HB2	1:B:151:GLU:HB2	1.90	0.53
1:A:54:GLN:HG2	1:A:58:ASN:HD21	1.73	0.53
1:A:67:PRO:HG3	1:A:89:THR:HG23	1.90	0.53
1:B:122:SER:O	1:B:126:GLN:HG3	2.10	0.52
1:B:52:PHE:CE1	1:B:61:ILE:HD13	2.44	0.52
1:B:94:MSE:HG3	1:B:123:TYR:HB3	1.91	0.52
1:B:269:ASP:HB2	3:B:1198:HOH:O	2.09	0.52
1:B:94:MSE:HE2	1:B:127:ALA:HB2	1.90	0.51
1:A:229:MSE:HB2	1:B:173:MSE:HE1	1.91	0.51
1:A:177:MSE:HG2	3:A:1140:HOH:O	2.09	0.51
1:B:62:ILE:HD12	1:B:83:ARG:HB2	1.92	0.51
1:A:204:ASN:HB2	3:A:1088:HOH:O	2.11	0.51
1:B:67:PRO:HG3	1:B:89:THR:HG23	1.92	0.51
1:A:22:ASP:HB3	1:A:25:THR:HB	1.91	0.51
1:B:56:LEU:HD11	1:B:63:PRO:HG3	1.93	0.51
1:B:159:LEU:C	1:B:160:ILE:HD12	2.32	0.50
1:B:157:ASN:N	1:B:157:ASN:ND2	2.58	0.50
1:B:177:MSE:O	1:B:177:MSE:HE3	2.12	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:214:LYS:HE3	3:A:1142:HOH:O	2.11	0.50
1:B:259:ILE:HG23	1:B:260:SER:N	2.26	0.50
1:B:81:ALA:HB2	1:B:111:LEU:CD2	2.41	0.50
1:A:249:VAL:HA	1:A:253:ASN:ND2	2.27	0.49
1:A:186:ASN:HD22	1:A:191:ARG:HD2	1.77	0.49
1:A:9:LYS:CE	1:A:31:GLU:HG2	2.41	0.49
1:B:249:VAL:HA	1:B:253:ASN:HD22	1.77	0.49
1:A:292:GLU:O	1:B:150:LYS:HG2	2.13	0.49
1:A:157:ASN:ND2	1:A:157:ASN:H	2.10	0.49
1:B:150:LYS:HB2	1:B:217:ARG:HH21	1.78	0.48
1:A:67:PRO:CG	1:A:89:THR:HG23	2.43	0.48
1:B:139:ASN:OD1	1:B:259:ILE:HG13	2.14	0.48
1:B:325:VAL:O	1:B:329:ILE:HG13	2.14	0.48
1:A:118:MSE:HE3	3:A:1056:HOH:O	2.14	0.48
1:A:300:ILE:CD1	1:B:149:ALA:HB2	2.44	0.47
1:A:173:MSE:HE1	1:B:225:LYS:O	2.15	0.47
1:B:150:LYS:HE3	1:B:217:ARG:HH21	1.80	0.47
1:A:61:ILE:HD12	1:A:61:ILE:N	2.30	0.47
1:A:154:ILE:HG22	1:A:155:GLY:N	2.30	0.47
1:B:160:ILE:HD12	1:B:160:ILE:N	2.29	0.46
1:A:221:MSE:HE3	1:B:221:MSE:SE	2.65	0.46
1:B:141:ILE:O	1:B:143:VAL:HG13	2.16	0.46
1:A:154:ILE:HD11	1:B:288:VAL:CG2	2.44	0.46
1:B:94:MSE:CE	1:B:98:ILE:HD11	2.46	0.46
1:B:210:GLY:O	1:B:214:LYS:HG2	2.16	0.46
1:B:219:LEU:HD12	1:B:220:PRO:HD2	1.98	0.45
1:B:249:VAL:HA	1:B:253:ASN:ND2	2.30	0.45
1:B:79:TRP:HB3	1:B:93:MSE:CE	2.46	0.45
1:A:193:LEU:CD2	1:A:209:PRO:HG3	2.46	0.45
1:B:169:THR:OG1	1:B:305:HIS:HE1	2.00	0.45
1:A:113:ASN:HA	1:A:258:GLY:HA2	1.98	0.45
1:A:214:LYS:HD2	1:A:214:LYS:C	2.35	0.45
1:A:85:ILE:HD12	1:A:88:LEU:HD12	1.99	0.45
1:A:8:LEU:HD22	1:A:12:PHE:CZ	2.51	0.45
1:A:153:ARG:HD2	1:B:287:TYR:CE1	2.52	0.45
1:A:186:ASN:ND2	1:A:191:ARG:HD2	2.31	0.44
1:B:112:ARG:HB3	1:B:241:SER:HB3	1.99	0.44
1:A:159:LEU:C	1:A:160:ILE:HD12	2.38	0.44
1:A:157:ASN:N	1:A:157:ASN:ND2	2.63	0.44
1:A:8:LEU:HD22	1:A:12:PHE:CE1	2.52	0.44
1:B:209:PRO:O	1:B:213:GLU:HG3	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:187:ARG:HH11	1:B:187:ARG:HG2	1.83	0.44
1:A:160:ILE:HD12	1:A:160:ILE:N	2.33	0.44
1:B:4:THR:HG22	1:B:7:GLN:CB	2.46	0.44
1:B:139:ASN:HA	1:B:157:ASN:ND2	2.33	0.43
1:A:198:GLY:O	1:A:205:LEU:HA	2.18	0.43
1:B:150:LYS:CB	1:B:217:ARG:HH21	2.31	0.43
1:B:215:ASN:N	1:B:215:ASN:ND2	2.66	0.43
1:B:49:PHE:CZ	1:B:53:ILE:HD11	2.52	0.43
1:A:150:LYS:HE3	1:B:293:ARG:O	2.19	0.43
1:A:9:LYS:NZ	1:A:31:GLU:HG2	2.34	0.43
1:B:208:GLU:O	1:B:212:ILE:HG13	2.18	0.43
1:A:63:PRO:HG2	3:A:1258:HOH:O	2.17	0.43
1:A:93:MSE:HE2	1:A:93:MSE:HA	2.00	0.43
1:A:301:ARG:NH1	1:A:301:ARG:HG2	2.33	0.43
1:A:31:GLU:HG3	3:A:1303:HOH:O	2.19	0.43
1:B:108:LEU:HD23	1:B:108:LEU:C	2.39	0.43
1:A:60:ASP:HB2	1:A:61:ILE:HD12	2.00	0.42
1:A:324:SER:O	1:A:328:LYS:HD3	2.19	0.42
1:A:300:ILE:HD13	1:B:149:ALA:HB2	2.02	0.42
1:A:85:ILE:HG13	3:A:1187:HOH:O	2.18	0.42
1:B:93:MSE:HE3	1:B:111:LEU:HG	2.02	0.42
1:A:1:MSE:HA	1:A:1:MSE:HE2	2.01	0.42
1:B:277:ARG:O	1:B:281:LEU:HB2	2.20	0.42
1:A:204:ASN:ND2	1:A:205:LEU:N	2.64	0.41
1:A:154:ILE:HD12	1:A:231:ILE:HG12	2.02	0.41
1:B:273:ASP:CG	1:B:275:PRO:HD2	2.41	0.41
1:A:311:LEU:O	1:A:315:ARG:HG3	2.20	0.41
1:B:44:HIS:HD2	3:B:1109:HOH:O	2.04	0.41
1:B:181:GLY:O	1:B:185:VAL:HG23	2.19	0.41
1:B:221:MSE:O	1:B:226:GLY:HA3	2.20	0.41
1:A:106:ILE:HG22	1:B:76:ILE:HD11	2.02	0.41
1:A:160:ILE:HG13	1:A:172:ASP:HA	2.02	0.41
1:A:167:PRO:HG3	3:A:1153:HOH:O	2.20	0.41
1:B:180:TYR:CE2	1:B:218:ILE:HD13	2.56	0.41
1:B:48:ARG:HD3	1:B:51:ARG:HB3	2.02	0.41
1:A:146:PRO:HD3	1:A:154:ILE:HG12	2.02	0.41
1:B:215:ASN:HD22	1:B:216:ARG:N	2.18	0.41
1:B:71:THR:HG22	1:B:73:LEU:HG	2.02	0.41
1:A:38:GLU:HG2	1:A:311:LEU:HD12	2.01	0.41
1:A:12:PHE:CE1	1:A:33:PHE:HB3	2.56	0.40
1:A:283:ARG:NH1	1:B:248:GLU:HG3	2.35	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:298:GLN:HA	1:A:298:GLN:OE1	2.20	0.40
1:B:192:GLN:HE22	1:B:208:GLU:HG2	1.85	0.40
1:B:188:LEU:HB3	1:B:322:ASP:OD2	2.21	0.40
1:A:186:ASN:HD22	1:A:191:ARG:CG	2.34	0.40
1:A:147:TRP:CD1	1:A:224:TRP:HB3	2.57	0.40
1:B:146:PRO:HD3	1:B:153:ARG:HB2	2.02	0.40
1:B:329:ILE:O	1:B:332:LEU:HB2	2.21	0.40
1:B:12:PHE:CE1	1:B:33:PHE:HB3	2.57	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	330/333 (99%)	312 (94%)	16 (5%)	2 (1%)	30	14
1	B	331/333 (99%)	316 (96%)	14 (4%)	1 (0%)	46	29
All	All	661/666 (99%)	628 (95%)	30 (4%)	3 (0%)	34	17

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	201	ASP
1	A	118	MSE
1	B	190	GLY

5.3.2 Protein sidechains [i](#)

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	271/256 (106%)	264 (97%)	7 (3%)	54	37
1	B	271/256 (106%)	257 (95%)	14 (5%)	29	12
All	All	542/512 (106%)	521 (96%)	21 (4%)	39	21

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MSE
1	A	117	TRP
1	A	118	MSE
1	A	144	MSE
1	A	204	ASN
1	A	214	LYS
1	A	281	LEU
1	B	2	LYS
1	B	88	LEU
1	B	111	LEU
1	B	117	TRP
1	B	193	LEU
1	B	201	ASP
1	B	204	ASN
1	B	215	ASN
1	B	259	ILE
1	B	273	ASP
1	B	281	LEU
1	B	301	ARG
1	B	305	HIS
1	B	316	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	44	HIS
1	A	54	GLN
1	A	58	ASN
1	A	78	GLN
1	A	82	GLN
1	A	104	HIS

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Mol	Chain	Res	Type
1	A	113	ASN
1	A	157	ASN
1	A	186	ASN
1	A	192	GLN
1	A	204	ASN
1	A	253	ASN
1	A	314	ASN
1	B	44	HIS
1	B	157	ASN
1	B	192	GLN
1	B	215	ASN
1	B	251	GLN
1	B	305	HIS

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](#)

4 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	A	1001	-	4,4,4	0.15	0	6,6,6	0.14	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	1002	-	4,4,4	0.21	0	6,6,6	0.12	0
2	SO4	B	1003	-	4,4,4	0.22	0	6,6,6	0.09	0
2	SO4	B	1004	-	4,4,4	0.14	0	6,6,6	0.06	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	SO4	A	1001	-	-	0/0/0/0	0/0/0/0
2	SO4	A	1002	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1003	-	-	0/0/0/0	0/0/0/0
2	SO4	B	1004	-	-	0/0/0/0	0/0/0/0

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data ⓘ

6.1 Protein, DNA and RNA chains ⓘ

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	317/333 (95%)	0.09	8 (2%) 61 56	10, 22, 43, 59	0
1	B	318/333 (95%)	0.20	14 (4%) 38 32	12, 23, 51, 68	0
All	All	635/666 (95%)	0.15	22 (3%) 48 42	10, 22, 46, 68	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	202	GLU	5.4
1	B	201	ASP	4.7
1	A	203	GLY	3.9
1	B	202	GLU	3.8
1	A	201	ASP	3.3
1	A	64	ASP	3.3
1	B	217	ARG	3.2
1	A	204	ASN	3.1
1	B	204	ASN	2.9
1	B	203	GLY	2.9
1	B	214	LYS	2.9
1	B	200	ASP	2.8
1	A	63	PRO	2.7
1	B	198	GLY	2.7
1	B	199	PHE	2.6
1	B	2	LYS	2.5
1	B	205	LEU	2.5
1	B	212	ILE	2.4
1	B	211	VAL	2.3
1	A	214	LYS	2.3
1	A	159	LEU	2.2
1	B	159	LEU	2.1

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, 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	LLDF	B-factors(\AA^2)	Q<0.9
2	SO4	A	1002	5/5	0.98	0.07	-1.36	17,19,21,22	0
2	SO4	B	1004	5/5	0.98	0.09	-1.65	22,22,24,25	0
2	SO4	A	1001	5/5	0.98	0.07	-1.83	24,26,26,27	0
2	SO4	B	1003	5/5	0.97	0.08	-1.96	28,29,30,31	0

6.5 Other polymers [i](#)

There are no such residues in this entry.