



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 01:54 PM UTC

PDB ID : 2X0X / pdb_00002x0x
Title : Ribonucleotide reductase R1 subunit of E. coli to 2.3 Å resolution
Authors : Yokoyama, K.; Uhlin, U.; Stubbe, J.
Deposited on : 2009-12-18
Resolution : 2.30 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

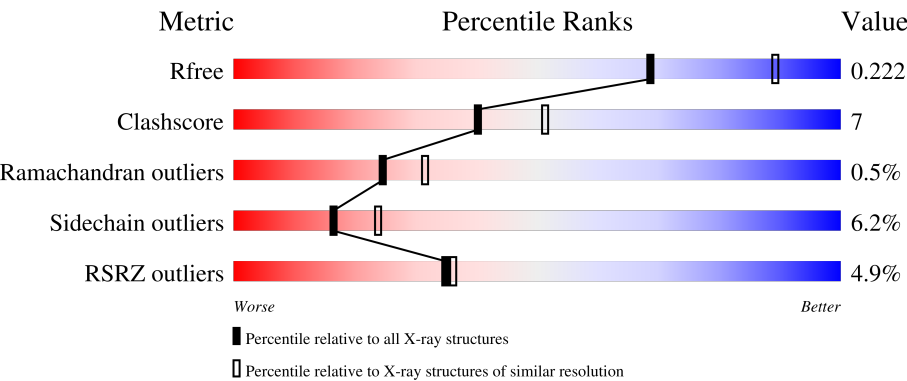
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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	761	4% 79% 14% • •
1	B	761	4% 81% 13% • •
1	C	761	3% 81% 12% • •
2	D	20	45% 70% 10% 20%
2	E	20	35% 70% 5% 5% 20%

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Mol	Chain	Length	Quality of chain
2	F	20	25% 60% 10% 10% 20%
2	P	20	5% 5% 10% 85%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 18869 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 RIBONUCLEOSIDE-DIPHOSPHATE REDUCTASE 1 SUBUNIT ALPHA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	728	Total	C	N	O	S	0	0	0
			5804	3688	995	1097	24			
1	B	728	Total	C	N	O	S	0	0	0
			5804	3688	995	1097	24			
1	C	728	Total	C	N	O	S	0	0	0
			5804	3688	995	1097	24			

- Molecule 2 is a protein called RIBONUCLEOSIDE-DIPHOSPHATE REDUCTASE 1 SUBUNIT BETA.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
2	D	16	Total	C	N	O	0	0	0
			129	77	19	33			
2	E	16	Total	C	N	O	0	0	0
			129	77	19	33			
2	F	16	Total	C	N	O	0	0	0
			129	77	19	33			
2	P	3	Total	C	N	O	0	0	0
			27	20	3	4			

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

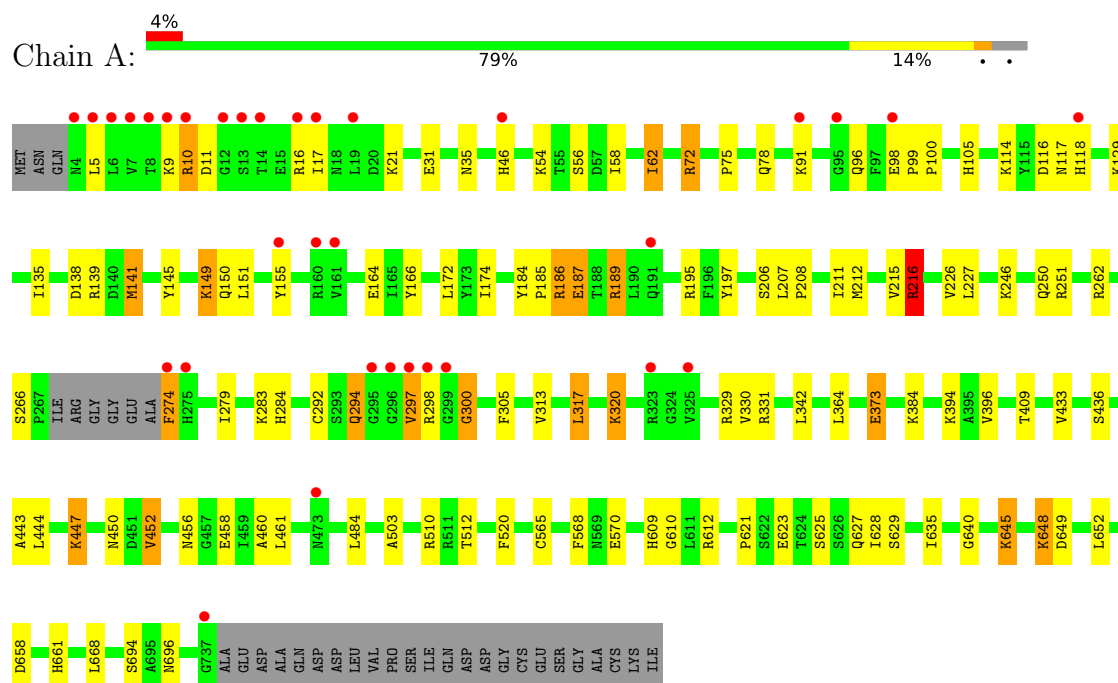
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	304	Total	O	0	0
			304	304		
4	B	306	Total	O	0	0
			306	306		
4	C	411	Total	O	0	0
			411	411		
4	D	1	Total	O	0	0
			1	1		
4	E	3	Total	O	0	0
			3	3		
4	F	2	Total	O	0	0
			2	2		
4	P	1	Total	O	0	0
			1	1		

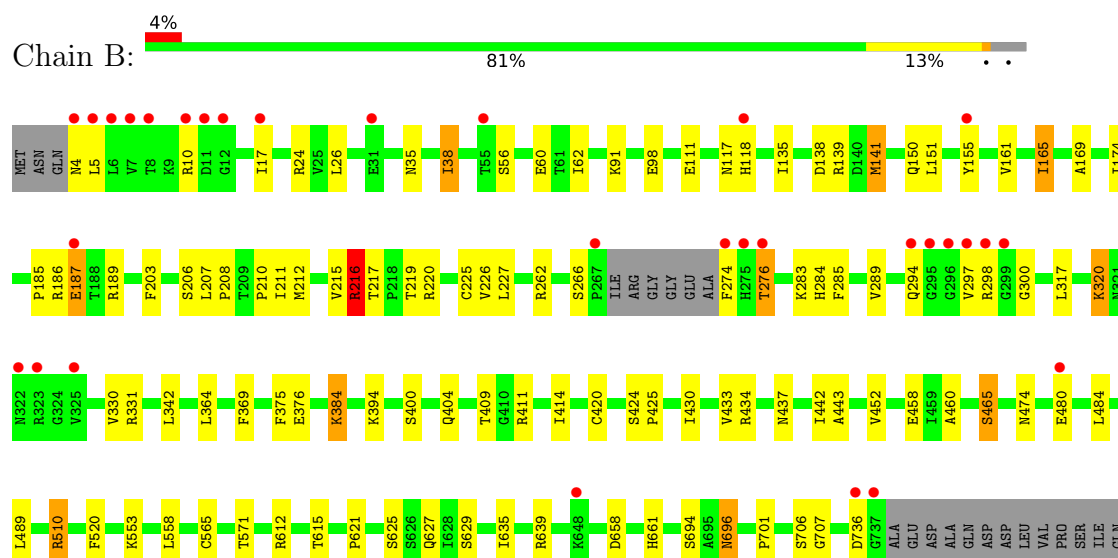
3 Residue-property plots

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: RIBONUCLEOSIDE-DIPHOSPHATE REDUCTASE 1 SUBUNIT ALPHA



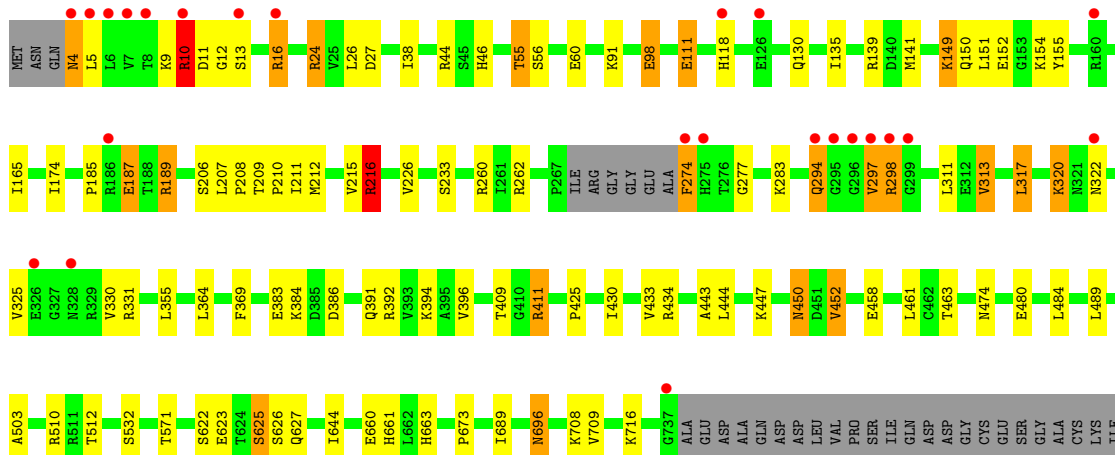
• Molecule 1: RIBONUCLEOSIDE-DIPHOSPHATE REDUCTASE 1 SUBUNIT ALPHA



ASP
ASP
GLY
CYS
GLU
SER
GLY
ALA
CYS
LYS
ILE

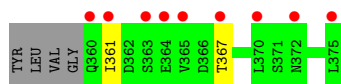
• Molecule 1: RIBONUCLEOSIDE-DIPHOSPHATE REDUCTASE 1 SUBUNIT ALPHA

Chain C: 



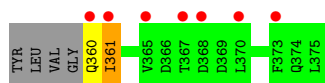
• Molecule 2: RIBONUCLEOSIDE-DIPHOSPHATE REDUCTASE 1 SUBUNIT BETA

Chain D: 



• Molecule 2: RIBONUCLEOSIDE-DIPHOSPHATE REDUCTASE 1 SUBUNIT BETA

Chain E: 



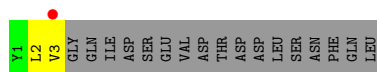
• Molecule 2: RIBONUCLEOSIDE-DIPHOSPHATE REDUCTASE 1 SUBUNIT BETA

Chain F: 



• Molecule 2: RIBONUCLEOSIDE-DIPHOSPHATE REDUCTASE 1 SUBUNIT BETA

Chain P: 



4 Data and refinement statistics

Property	Value	Source
Space group	H 3 2	Depositor
Cell constants a, b, c, α , β , γ	223.88Å 223.88Å 336.47Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	169.03 – 2.30 167.99 – 2.30	Depositor EDS
% Data completeness (in resolution range)	100.0 (169.03-2.30) 100.0 (167.99-2.30)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.35 (at 2.29Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.183 , 0.224 0.182 , 0.222	Depositor DCC
R_{free} test set	7167 reflections (5.01%)	wwPDB-VP
Wilson B-factor (Å ²)	34.6	Xtriage
Anisotropy	0.120	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.36 , 39.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	18869	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.87% 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	0.67	1/5931 (0.0%)	0.90	1/8033 (0.0%)
1	B	0.67	0/5931	0.90	3/8033 (0.0%)
1	C	0.77	1/5931 (0.0%)	0.95	8/8033 (0.1%)
2	D	0.60	0/129	0.87	0/173
2	E	0.62	0/129	0.94	0/173
2	F	0.62	0/129	0.85	0/173
2	P	0.71	0/27	0.57	0/36
All	All	0.70	2/18207 (0.0%)	0.91	12/24654 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	689	ILE	CA-CB	6.00	1.61	1.54
1	A	279	ILE	CA-CB	5.54	1.57	1.54

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	452	VAL	CB-CA-C	-6.92	101.95	112.05
1	B	38	ILE	CB-CA-C	-6.28	103.84	111.88
1	C	430	ILE	CB-CA-C	-6.10	105.65	111.44
1	C	277	GLY	N-CA-C	5.66	118.94	111.70
1	A	452	VAL	CB-CA-C	-5.57	104.62	112.14
1	B	430	ILE	CB-CA-C	-5.51	106.21	111.44
1	C	216	ARG	NE-CZ-NH2	-5.37	114.37	119.20
1	C	313	VAL	N-CA-C	5.34	116.70	110.62
1	C	154	LYS	N-CA-C	5.34	121.34	114.56
1	C	216	ARG	CG-CD-NE	-5.26	100.42	112.00
1	B	169	ALA	N-CA-C	5.07	116.81	111.28
1	C	450	ASN	CB-CA-C	-5.02	100.91	109.65

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	5804	0	5728	88	0
1	B	5804	0	5728	60	0
1	C	5804	0	5728	83	0
2	D	129	0	111	0	0
2	E	129	0	111	1	0
2	F	129	0	111	3	0
2	P	27	0	31	5	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	1	0
4	A	304	0	0	18	0
4	B	306	0	0	12	1
4	C	411	0	0	21	1
4	D	1	0	0	0	0
4	E	3	0	0	0	0
4	F	2	0	0	1	0
4	P	1	0	0	2	0
All	All	18869	0	17548	233	1

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 7.

All (233) 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:342:LEU:HG	4:A:2133:HOH:O	1.38	1.21
1:C:189:ARG:NH1	4:C:2117:HOH:O	1.87	1.07
1:A:186:ARG:HG3	1:A:186:ARG:HH11	1.25	0.99
1:B:274:PHE:HA	4:B:2117:HOH:O	1.63	0.96
1:C:450:ASN:HB2	4:C:2242:HOH:O	1.67	0.92
2:P:3:VAL:HA	4:P:2001:HOH:O	1.73	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:155:TYR:CD1	4:C:2095:HOH:O	2.28	0.87
1:C:480:GLU:HB3	4:C:2120:HOH:O	1.74	0.85
1:A:262:ARG:CG	1:A:274:PHE:HB2	2.06	0.85
1:C:155:TYR:HD1	4:C:2095:HOH:O	1.60	0.83
1:B:262:ARG:HD3	1:B:274:PHE:HB3	1.62	0.82
1:B:56:SER:O	1:B:60:GLU:HG2	1.79	0.81
1:B:215:VAL:O	1:B:216:ARG:HB3	1.81	0.80
1:B:38:ILE:HG12	4:B:2005:HOH:O	1.82	0.80
1:A:10:ARG:H	1:A:10:ARG:HD3	1.46	0.80
1:C:44:ARG:HD3	4:C:2025:HOH:O	1.81	0.79
1:C:262:ARG:HG2	1:C:274:PHE:HB2	1.65	0.79
1:A:212:MET:O	1:A:216:ARG:NH2	2.14	0.79
1:A:450:ASN:HB2	4:A:2175:HOH:O	1.84	0.77
1:C:215:VAL:O	1:C:216:ARG:HB3	1.84	0.77
1:C:298:ARG:HG2	1:C:298:ARG:HH11	1.48	0.77
1:B:185:PRO:HB2	1:B:187:GLU:HG2	1.67	0.76
1:A:9:LYS:HG3	1:A:10:ARG:NH1	2.00	0.76
1:C:262:ARG:CG	1:C:274:PHE:HB2	2.17	0.75
1:C:208:PRO:HD2	1:C:211:ILE:HD12	1.69	0.74
1:A:215:VAL:O	1:A:216:ARG:HB3	1.87	0.74
1:B:208:PRO:HD2	1:B:211:ILE:HD12	1.71	0.73
1:A:250:GLN:NE2	4:A:2105:HOH:O	2.22	0.73
1:C:189:ARG:NH2	4:C:2116:HOH:O	2.17	0.72
1:C:260:ARG:HG2	1:C:260:ARG:HH11	1.56	0.71
1:C:24:ARG:HH11	1:C:24:ARG:CG	2.03	0.71
1:A:648:LYS:H	1:A:648:LYS:HE2	1.56	0.69
1:A:186:ARG:HG3	1:A:186:ARG:NH1	2.01	0.69
1:C:189:ARG:NE	4:C:2116:HOH:O	2.09	0.68
1:A:262:ARG:HG3	1:A:274:PHE:HB2	1.75	0.67
1:C:262:ARG:HG2	1:C:274:PHE:CB	2.25	0.67
1:C:283:LYS:HG3	1:C:330:VAL:HG22	1.77	0.66
1:A:195:ARG:HD3	4:A:2063:HOH:O	1.97	0.65
1:C:24:ARG:HH11	1:C:24:ARG:HG3	1.61	0.65
1:C:450:ASN:CB	4:C:2242:HOH:O	2.36	0.65
1:A:262:ARG:HG2	1:A:274:PHE:HB2	1.78	0.65
1:C:622:SER:HG	1:C:626:SER:HG	1.45	0.65
1:C:209:THR:HG23	4:C:2095:HOH:O	1.98	0.64
1:A:648:LYS:HD2	1:A:649:ASP:H	1.63	0.63
1:A:54:LYS:HE2	1:A:56:SER:H	1.62	0.63
1:B:151:LEU:HA	1:B:155:TYR:HB2	1.81	0.63
1:B:212:MET:O	1:B:216:ARG:NH2	2.24	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:396:VAL:HG23	4:A:2137:HOH:O	1.99	0.62
1:A:118:HIS:CE1	4:A:2035:HOH:O	2.52	0.62
1:A:292:CYS:HA	1:B:276:THR:HG21	1.80	0.62
1:C:625:SER:OG	3:C:1738:SO4:O4	2.15	0.62
1:A:320:LYS:HA	1:A:331:ARG:HG2	1.81	0.62
1:C:27:ASP:HA	1:C:38:ILE:HD11	1.80	0.62
1:A:262:ARG:HG2	1:A:274:PHE:CB	2.31	0.60
1:C:396:VAL:HG23	4:C:2197:HOH:O	2.01	0.60
1:A:433:VAL:HG11	1:A:443:ALA:HB1	1.82	0.60
1:A:313:VAL:HG22	1:A:317:LEU:HD22	1.83	0.59
1:B:24:ARG:HD2	4:B:2004:HOH:O	2.03	0.59
1:A:16:ARG:HG3	4:A:2001:HOH:O	2.03	0.59
1:A:207:LEU:HD12	1:A:212:MET:HE3	1.84	0.59
1:A:91:LYS:NZ	4:A:2026:HOH:O	2.34	0.58
1:A:151:LEU:HA	1:A:155:TYR:HB2	1.84	0.58
1:C:10:ARG:HH11	1:C:10:ARG:HB3	1.68	0.58
1:A:186:ARG:HH11	1:A:186:ARG:CG	2.09	0.58
1:B:187:GLU:HB2	4:B:2077:HOH:O	2.03	0.58
1:B:639:ARG:HD3	4:B:2260:HOH:O	2.02	0.58
1:A:75:PRO:O	1:A:78:GLN:HB2	2.04	0.58
4:C:2196:HOH:O	2:F:367:THR:HG22	2.02	0.57
1:C:260:ARG:HG2	1:C:260:ARG:NH1	2.19	0.57
1:A:207:LEU:CD1	1:A:212:MET:HE3	2.34	0.57
1:A:510:ARG:NH2	1:A:570:GLU:OE1	2.33	0.57
1:C:189:ARG:HH11	1:C:189:ARG:CG	2.17	0.57
1:B:207:LEU:HB2	1:B:212:MET:HE3	1.85	0.57
1:C:433:VAL:HG11	1:C:443:ALA:HB1	1.85	0.57
1:C:207:LEU:HB2	1:C:212:MET:HE3	1.87	0.57
1:A:458:GLU:OE2	1:A:510:ARG:HD2	2.04	0.56
1:A:444:LEU:HD22	1:A:512:THR:HG21	1.86	0.56
1:A:329:ARG:HG3	1:A:331:ARG:NH1	2.21	0.56
1:C:155:TYR:CZ	1:C:212:MET:HG3	2.41	0.56
1:C:260:ARG:HH21	1:C:434:ARG:HH22	1.54	0.56
1:C:274:PHE:N	1:C:274:PHE:CD2	2.74	0.56
1:B:465:SER:HB2	1:B:489:LEU:HD11	1.88	0.55
1:A:300:GLY:HA2	4:A:2106:HOH:O	2.06	0.55
1:C:111:GLU:HG3	2:P:2:LEU:HB2	1.88	0.55
1:A:207:LEU:HB2	1:A:212:MET:HE3	1.88	0.55
1:C:189:ARG:HH11	1:C:189:ARG:HG2	1.72	0.55
1:A:373:GLU:HG2	4:A:2150:HOH:O	2.06	0.55
1:A:262:ARG:HD3	1:A:266:SER:HB2	1.88	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:409:THR:CG2	4:A:2156:HOH:O	2.54	0.55
1:A:138:ASP:HA	1:A:141:MET:HE2	1.88	0.55
1:A:227:LEU:HB2	1:A:460:ALA:HB3	1.89	0.55
1:C:24:ARG:HD2	4:C:2008:HOH:O	2.06	0.54
1:A:658:ASP:OD1	1:A:661:HIS:HD2	1.89	0.54
1:A:294:GLN:HB2	1:A:298:ARG:HH11	1.73	0.53
2:F:361:ILE:HD11	4:F:2001:HOH:O	2.08	0.53
1:C:313:VAL:HG22	1:C:317:LEU:HD22	1.89	0.53
1:A:58:ILE:O	1:A:62:ILE:HG23	2.09	0.53
1:C:24:ARG:CG	1:C:24:ARG:NH1	2.69	0.53
1:B:135:ILE:HD11	1:B:174:ILE:HG21	1.91	0.52
1:C:458:GLU:OE2	1:C:510:ARG:HD2	2.08	0.52
1:B:696:ASN:N	1:B:696:ASN:HD22	2.08	0.52
1:B:283:LYS:HG3	1:B:330:VAL:HG22	1.92	0.52
2:P:2:LEU:O	2:P:3:VAL:C	2.52	0.52
1:A:447:LYS:HD2	1:A:456:ASN:O	2.10	0.51
1:B:520:PHE:HB3	1:B:635:ILE:HA	1.92	0.51
1:C:150:GLN:HE21	1:C:627:GLN:CD	2.18	0.51
1:A:116:ASP:CG	1:A:118:HIS:HD2	2.17	0.51
1:A:208:PRO:HD2	1:A:211:ILE:HD12	1.91	0.51
1:C:155:TYR:CE2	1:C:212:MET:SD	3.03	0.51
1:C:709:VAL:O	2:F:361:ILE:HB	2.11	0.51
1:B:117:ASN:HB2	4:B:2043:HOH:O	2.09	0.51
1:A:305:PHE:CZ	1:A:436:SER:HB3	2.46	0.50
1:B:658:ASP:OD1	1:B:661:HIS:HD2	1.95	0.50
1:A:117:ASN:HB2	4:A:2034:HOH:O	2.11	0.50
1:C:409:THR:O	1:C:411:ARG:HG2	2.12	0.50
1:A:186:ARG:NH1	1:A:186:ARG:CG	2.71	0.49
1:B:701:PRO:O	1:B:707:GLY:HA2	2.12	0.49
1:C:320:LYS:HA	1:C:331:ARG:HG2	1.95	0.49
1:C:189:ARG:NH1	1:C:189:ARG:HG2	2.28	0.49
1:B:342:LEU:HD23	1:B:375:PHE:HE2	1.78	0.49
1:B:420:CYS:O	1:B:424:SER:HB2	2.13	0.49
1:C:383:GLU:O	1:C:392:ARG:NH2	2.46	0.49
1:C:320:LYS:HE2	1:C:411:ARG:HB2	1.95	0.48
1:B:10:ARG:NH1	1:B:91:LYS:HD3	2.27	0.48
1:B:207:LEU:HD22	1:B:211:ILE:HG21	1.95	0.48
1:A:409:THR:HG23	4:A:2156:HOH:O	2.14	0.48
1:A:648:LYS:HD2	1:A:649:ASP:N	2.28	0.48
1:C:56:SER:O	1:C:60:GLU:HG2	2.14	0.48
1:C:10:ARG:CZ	1:C:55:THR:HG21	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:294:GLN:HB3	4:C:2173:HOH:O	2.13	0.48
1:A:164:GLU:HG3	4:A:2028:HOH:O	2.13	0.48
1:B:404:GLN:HG3	2:E:361:ILE:HD11	1.96	0.48
1:A:185:PRO:HB2	1:A:187:GLU:OE2	2.14	0.47
1:A:9:LYS:HG3	1:A:10:ARG:HH11	1.77	0.47
1:A:283:LYS:HD3	4:B:2040:HOH:O	2.14	0.47
1:C:274:PHE:HB3	4:C:2168:HOH:O	2.14	0.47
1:B:474:ASN:HB3	4:B:2189:HOH:O	2.13	0.47
1:B:736:ASP:OD1	1:B:736:ASP:O	2.33	0.47
1:C:298:ARG:HG2	1:C:298:ARG:NH1	2.25	0.47
1:C:10:ARG:HB3	1:C:10:ARG:NH1	2.30	0.47
1:C:135:ILE:HD11	1:C:174:ILE:HG21	1.97	0.47
1:B:558:LEU:HD23	1:B:612:ARG:HG2	1.97	0.47
1:C:298:ARG:HH11	1:C:298:ARG:CG	2.24	0.46
1:B:433:VAL:HG11	1:B:443:ALA:HB1	1.96	0.46
1:A:262:ARG:CG	1:A:274:PHE:CB	2.83	0.46
1:B:165:ILE:HD13	4:B:2059:HOH:O	2.15	0.46
1:C:461:LEU:HD11	1:C:503:ALA:HB1	1.97	0.46
1:B:227:LEU:HB2	1:B:460:ALA:HB3	1.98	0.46
1:C:16:ARG:NE	1:C:16:ARG:H	2.13	0.46
1:C:663:HIS:HD2	4:C:2356:HOH:O	1.98	0.46
1:A:10:ARG:HD3	1:A:10:ARG:N	2.22	0.46
1:A:145:TYR:CZ	1:A:149:LYS:HD3	2.51	0.46
1:B:384:LYS:HA	1:B:384:LYS:HE3	1.96	0.46
1:B:409:THR:O	1:B:411:ARG:HG2	2.16	0.45
1:C:151:LEU:HA	1:C:155:TYR:HB2	1.98	0.45
1:A:609:HIS:HD2	4:C:2058:HOH:O	1.99	0.45
1:A:135:ILE:HD11	1:A:174:ILE:HG21	1.99	0.45
1:B:118:HIS:HB3	4:B:2044:HOH:O	2.17	0.45
1:C:660:GLU:OE2	1:C:661:HIS:NE2	2.50	0.45
1:A:100:PRO:HG2	1:A:105:HIS:HB2	1.99	0.45
1:A:640:GLY:HA2	1:A:668:LEU:HD13	1.99	0.45
1:C:212:MET:O	1:C:216:ARG:NH2	2.39	0.45
1:A:568:PHE:CE2	1:A:610:GLY:HA2	2.52	0.45
1:C:149:LYS:HE2	1:C:152:GLU:OE1	2.17	0.45
1:B:285:PHE:O	1:B:289:VAL:HG23	2.17	0.44
1:C:9:LYS:O	1:C:11:ASP:N	2.49	0.44
1:B:26:LEU:HB3	1:B:38:ILE:HG23	2.00	0.44
1:A:184:TYR:O	1:A:189:ARG:HD3	2.18	0.44
1:B:203:PHE:HB3	1:B:629:SER:HB2	1.98	0.44
1:B:217:THR:OG1	1:B:219:THR:HG22	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:444:LEU:HD22	1:C:512:THR:HG21	1.98	0.44
1:C:532:SER:HB2	1:C:673:PRO:HD2	2.00	0.44
1:A:215:VAL:O	1:A:216:ARG:CB	2.62	0.44
1:A:150:GLN:HE21	1:A:627:GLN:CD	2.25	0.44
1:C:10:ARG:CD	1:C:10:ARG:H	2.30	0.44
1:A:114:LYS:HE2	1:A:166:TYR:CE2	2.53	0.43
1:A:294:GLN:HB2	1:A:298:ARG:NH1	2.32	0.43
1:A:645:LYS:HD3	1:A:652:LEU:HB2	2.00	0.43
1:C:46:HIS:HB3	4:C:2026:HOH:O	2.17	0.43
1:C:463:THR:HG22	1:C:489:LEU:HD22	1.99	0.43
1:A:207:LEU:HD12	1:A:212:MET:CE	2.48	0.43
2:P:3:VAL:CA	4:P:2001:HOH:O	2.50	0.43
1:A:520:PHE:HB3	1:A:635:ILE:HA	2.01	0.43
1:A:246:LYS:NZ	4:A:2101:HOH:O	2.52	0.43
1:B:342:LEU:HD23	1:B:375:PHE:CE2	2.53	0.43
1:B:400:SER:O	1:B:404:GLN:HB2	2.19	0.43
1:A:172:LEU:C	1:A:172:LEU:HD23	2.44	0.42
1:A:621:PRO:HD3	1:A:694:SER:OG	2.19	0.42
1:B:150:GLN:HE21	1:B:627:GLN:CD	2.26	0.42
1:A:461:LEU:HD11	1:A:503:ALA:HB1	2.01	0.42
1:C:208:PRO:HB2	1:C:210:PRO:HD2	2.00	0.42
1:A:409:THR:HG21	4:A:2156:HOH:O	2.18	0.42
1:A:139:ARG:NH2	1:A:197:TYR:CE2	2.87	0.42
1:A:292:CYS:CA	1:B:276:THR:HG21	2.49	0.42
1:B:437:ASN:HB2	4:B:2174:HOH:O	2.19	0.42
1:C:311:LEU:HA	1:C:355:LEU:HB3	2.01	0.42
1:A:373:GLU:HG2	1:A:373:GLU:H	1.65	0.42
1:A:98:GLU:HA	1:A:99:PRO:HD3	1.83	0.42
1:A:565:CYS:HB3	1:A:612:ARG:O	2.19	0.42
1:B:621:PRO:HD3	1:B:694:SER:OG	2.20	0.42
1:C:10:ARG:NH1	1:C:55:THR:HG21	2.35	0.42
1:C:26:LEU:HB3	1:C:38:ILE:HG12	2.01	0.42
1:B:262:ARG:HG3	1:B:266:SER:HB2	2.02	0.42
1:A:283:LYS:HG3	1:A:330:VAL:HG22	2.01	0.42
1:A:284:HIS:CE1	1:B:284:HIS:CE1	3.08	0.42
1:C:369:PHE:CD2	1:C:434:ARG:HD2	2.54	0.42
1:C:708:LYS:HE2	1:C:708:LYS:HB3	1.74	0.42
1:A:450:ASN:HB2	4:A:2182:HOH:O	2.19	0.42
1:B:553:LYS:HE3	4:B:2226:HOH:O	2.20	0.41
1:B:320:LYS:HA	1:B:331:ARG:HG2	2.01	0.41
1:C:98:GLU:HG2	4:C:2054:HOH:O	2.19	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:425:PRO:O	1:B:571:THR:HB	2.21	0.41
1:B:565:CYS:HB3	1:B:612:ARG:O	2.21	0.41
1:C:425:PRO:O	1:C:571:THR:HB	2.21	0.41
1:A:72:ARG:H	1:A:72:ARG:HG2	1.52	0.41
1:B:225:CYS:SG	1:B:442:ILE:HG13	2.61	0.41
1:C:189:ARG:HH11	1:C:189:ARG:HB3	1.85	0.41
1:C:447:LYS:HE3	4:C:2254:HOH:O	2.20	0.41
1:B:208:PRO:HB2	1:B:210:PRO:HD2	2.02	0.41
1:C:111:GLU:CG	2:P:2:LEU:HB2	2.50	0.41
1:B:118:HIS:HD2	1:B:220:ARG:NH2	2.18	0.41
1:B:458:GLU:OE2	1:B:510:ARG:HD2	2.20	0.41
1:C:233:SER:HA	1:C:274:PHE:HZ	1.86	0.41
1:C:696:ASN:HD22	1:C:696:ASN:N	2.18	0.41
1:A:46:HIS:HE1	4:A:2005:HOH:O	2.04	0.40
1:B:138:ASP:O	1:B:141:MET:HB2	2.21	0.40
1:B:297:VAL:HG23	1:B:298:ARG:N	2.36	0.40
1:C:4:ASN:HD22	1:C:4:ASN:HA	1.79	0.40
1:B:369:PHE:CG	1:B:434:ARG:HG2	2.56	0.40
1:A:212:MET:C	1:A:216:ARG:HH22	2.19	0.40
1:C:185:PRO:HB2	1:C:187:GLU:HG2	2.02	0.40
1:B:425:PRO:HG2	1:B:615:THR:HG22	2.04	0.40
1:C:91:LYS:CG	4:C:2052:HOH:O	2.69	0.40

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:B:2006:HOH:O	4:C:2009:HOH:O[15_554]	2.16	0.04

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	724/761 (95%)	699 (96%)	21 (3%)	4 (1%)	21	27
1	B	724/761 (95%)	693 (96%)	28 (4%)	3 (0%)	30	38
1	C	724/761 (95%)	701 (97%)	18 (2%)	5 (1%)	18	23
2	D	14/20 (70%)	13 (93%)	1 (7%)	0	100	100
2	E	14/20 (70%)	13 (93%)	1 (7%)	0	100	100
2	F	14/20 (70%)	14 (100%)	0	0	100	100
2	P	1/20 (5%)	1 (100%)	0	0	100	100
All	All	2215/2363 (94%)	2134 (96%)	69 (3%)	12 (0%)	24	31

All (12) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	294	GLN
1	B	294	GLN
1	C	294	GLN
1	A	216	ARG
1	B	216	ARG
1	C	10	ARG
1	C	216	ARG
1	B	300	GLY
1	A	300	GLY
1	C	12	GLY
1	A	297	VAL
1	C	297	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	626/651 (96%)	588 (94%)	38 (6%)	17	24
1	B	626/651 (96%)	593 (95%)	33 (5%)	20	30
1	C	626/651 (96%)	585 (94%)	41 (6%)	15	22
2	D	16/19 (84%)	14 (88%)	2 (12%)	4	5

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	E	16/19 (84%)	14 (88%)	2 (12%)	4	5
2	F	16/19 (84%)	12 (75%)	4 (25%)	0	0
2	P	3/19 (16%)	3 (100%)	0	100	100
All	All	1929/2029 (95%)	1809 (94%)	120 (6%)	16	24

All (120) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	5	LEU
1	A	10	ARG
1	A	11	ASP
1	A	17	ILE
1	A	21	LYS
1	A	31	GLU
1	A	35	ASN
1	A	62	ILE
1	A	72	ARG
1	A	96	GLN
1	A	129	LYS
1	A	141	MET
1	A	149	LYS
1	A	186	ARG
1	A	187	GLU
1	A	189	ARG
1	A	206	SER
1	A	216	ARG
1	A	226	VAL
1	A	251	ARG
1	A	274	PHE
1	A	297	VAL
1	A	317	LEU
1	A	320	LYS
1	A	364	LEU
1	A	373	GLU
1	A	384	LYS
1	A	394	LYS
1	A	447	LYS
1	A	452	VAL
1	A	484	LEU
1	A	623	GLU
1	A	625	SER

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Mol	Chain	Res	Type
1	A	628	ILE
1	A	629	SER
1	A	645	LYS
1	A	648	LYS
1	A	696	ASN
1	B	4	ASN
1	B	5	LEU
1	B	17	ILE
1	B	35	ASN
1	B	62	ILE
1	B	98	GLU
1	B	111	GLU
1	B	139	ARG
1	B	141	MET
1	B	161	VAL
1	B	165	ILE
1	B	186	ARG
1	B	187	GLU
1	B	189	ARG
1	B	206	SER
1	B	216	ARG
1	B	226	VAL
1	B	276	THR
1	B	317	LEU
1	B	320	LYS
1	B	364	LEU
1	B	376	GLU
1	B	384	LYS
1	B	394	LYS
1	B	414	ILE
1	B	452	VAL
1	B	465	SER
1	B	480	GLU
1	B	484	LEU
1	B	510	ARG
1	B	625	SER
1	B	696	ASN
1	B	706	SER
1	C	4	ASN
1	C	5	LEU
1	C	10	ARG
1	C	13	SER

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Mol	Chain	Res	Type
1	C	16	ARG
1	C	24	ARG
1	C	55	THR
1	C	98	GLU
1	C	111	GLU
1	C	118	HIS
1	C	130	GLN
1	C	139	ARG
1	C	141	MET
1	C	149	LYS
1	C	165	ILE
1	C	187	GLU
1	C	189	ARG
1	C	206	SER
1	C	216	ARG
1	C	226	VAL
1	C	274	PHE
1	C	297	VAL
1	C	298	ARG
1	C	317	LEU
1	C	320	LYS
1	C	322	ASN
1	C	325	VAL
1	C	364	LEU
1	C	384	LYS
1	C	386	ASP
1	C	391	GLN
1	C	394	LYS
1	C	411	ARG
1	C	452	VAL
1	C	474	ASN
1	C	484	LEU
1	C	623	GLU
1	C	625	SER
1	C	644	ILE
1	C	696	ASN
1	C	716	LYS
2	D	361	ILE
2	D	367	THR
2	E	360	GLN
2	E	361	ILE
2	F	360	GLN

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Mol	Chain	Res	Type
2	F	361	ILE
2	F	367	THR
2	F	374	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	46	HIS
1	A	96	GLN
1	A	117	ASN
1	A	183	ASN
1	A	250	GLN
1	A	321	ASN
1	A	322	ASN
1	A	328	ASN
1	A	332	HIS
1	A	453	ASN
1	A	609	HIS
1	A	661	HIS
1	A	713	GLN
1	A	733	ASN
1	B	35	ASN
1	B	130	GLN
1	B	150	GLN
1	B	183	ASN
1	B	191	GLN
1	B	250	GLN
1	B	328	ASN
1	B	609	HIS
1	B	613	ASN
1	B	627	GLN
1	B	630	ASN
1	B	661	HIS
1	B	696	ASN
1	B	733	ASN
1	C	4	ASN
1	C	40	GLN
1	C	183	ASN
1	C	250	GLN
1	C	294	GLN
1	C	328	ASN
1	C	332	HIS

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Mol	Chain	Res	Type
1	C	456	ASN
1	C	474	ASN
1	C	630	ASN
1	C	633	ASN
1	C	663	HIS
1	C	696	ASN
1	C	698	ASN
1	C	733	ASN
2	D	360	GLN
2	E	360	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 oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

3 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	$\# Z > 2$
3	SO4	C	1738	-	4,4,4	0.30	0	6,6,6	0.26	0
3	SO4	A	1738	-	4,4,4	0.26	0	6,6,6	0.15	0
3	SO4	B	1738	-	4,4,4	0.31	0	6,6,6	0.23	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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1738	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, 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	728/761 (95%)	0.06	33 (4%) 38 40	22, 36, 58, 89	0
1	B	728/761 (95%)	0.17	31 (4%) 40 41	24, 37, 57, 84	0
1	C	728/761 (95%)	-0.20	24 (3%) 49 51	17, 28, 49, 75	0
2	D	16/20 (80%)	2.27	9 (56%) 0 0	70, 86, 91, 92	0
2	E	16/20 (80%)	2.09	7 (43%) 0 0	69, 85, 91, 92	0
2	F	16/20 (80%)	1.86	5 (31%) 1 1	57, 75, 82, 85	0
2	P	3/20 (15%)	1.36	1 (33%) 1 1	29, 29, 39, 45	0
All	All	2235/2363 (94%)	0.05	110 (4%) 35 36	17, 34, 63, 92	0

All (110) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	296	GLY	7.8
1	B	297	VAL	7.6
1	A	296	GLY	7.5
1	C	296	GLY	6.7
1	B	274	PHE	6.4
1	C	297	VAL	5.9
1	A	737	GLY	5.2
1	A	297	VAL	5.0
1	A	299	GLY	5.0
2	E	361	ILE	4.8
1	C	737	GLY	4.8
2	D	361	ILE	4.8
1	B	10	ARG	4.8
1	B	737	GLY	4.4
1	A	17	ILE	4.3
1	A	4	ASN	4.2
1	B	299	GLY	4.2

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Mol	Chain	Res	Type	RSRZ
2	F	361	ILE	4.2
1	A	274	PHE	4.1
1	C	274	PHE	4.1
1	B	11	ASP	4.1
1	A	5	LEU	4.1
1	C	295	GLY	3.9
1	A	6	LEU	3.8
1	B	5	LEU	3.5
1	A	118	HIS	3.5
1	B	6	LEU	3.4
1	A	7	VAL	3.4
1	B	8	THR	3.3
1	A	325	VAL	3.3
1	A	298	ARG	3.3
1	C	10	ARG	3.3
1	A	295	GLY	3.2
1	B	295	GLY	3.2
2	D	365	VAL	3.2
1	B	298	ARG	3.2
2	F	360	GLN	3.2
2	E	367	THR	3.2
1	B	736	ASP	3.1
2	P	3	VAL	3.1
1	B	4	ASN	3.1
1	B	17	ILE	3.1
1	C	294	GLN	3.1
1	A	161	VAL	3.0
1	C	299	GLY	3.0
2	D	372	ASN	3.0
1	B	118	HIS	3.0
1	B	294	GLN	3.0
1	B	275	HIS	2.9
1	A	10	ARG	2.9
2	E	365	VAL	2.9
1	A	13	SER	2.9
1	A	95	GLY	2.8
1	C	5	LEU	2.8
2	D	375	LEU	2.8
1	C	13	SER	2.7
1	C	4	ASN	2.7
1	C	160	ARG	2.7
1	B	267	PRO	2.7

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Mol	Chain	Res	Type	RSRZ
2	D	370	LEU	2.7
2	D	367	THR	2.7
2	F	367	THR	2.6
1	C	186	ARG	2.6
1	C	118	HIS	2.6
1	A	160	ARG	2.6
1	B	322	ASN	2.6
1	A	155	TYR	2.6
1	B	12	GLY	2.6
1	A	12	GLY	2.5
1	C	8	THR	2.5
1	C	326	GLU	2.5
2	D	364	GLU	2.5
2	E	373	PHE	2.5
1	B	7	VAL	2.5
1	A	19	LEU	2.5
2	E	360	GLN	2.5
1	B	55	THR	2.4
2	E	370	LEU	2.4
1	A	14	THR	2.4
2	D	363	SER	2.4
1	A	8	THR	2.4
1	B	480	GLU	2.4
1	C	126	GLU	2.4
1	C	328	ASN	2.3
1	C	298	ARG	2.3
1	B	323	ARG	2.3
2	D	360	GLN	2.3
1	A	323	ARG	2.3
1	B	276	THR	2.3
1	A	275	HIS	2.3
1	B	325	VAL	2.3
1	C	275	HIS	2.2
1	C	7	VAL	2.2
1	C	322	ASN	2.2
1	A	473	ASN	2.2
2	E	368	ASP	2.2
1	A	46	HIS	2.2
1	A	191	GLN	2.2
1	B	155	TYR	2.2
1	B	187	GLU	2.2
1	B	648	LYS	2.2

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Mol	Chain	Res	Type	RSRZ
1	C	6	LEU	2.1
1	C	16	ARG	2.1
1	A	91	LYS	2.1
1	A	16	ARG	2.1
2	F	375	LEU	2.1
1	A	9	LYS	2.0
2	F	365	VAL	2.0
1	A	98	GLU	2.0
1	B	31	GLU	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 oligosaccharides 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($\text{\AA}^2$)	Q<0.9
3	SO4	B	1738	5/5	0.75	0.22	123,124,124,124	0
3	SO4	C	1738	5/5	0.78	0.18	78,80,80,81	0
3	SO4	A	1738	5/5	0.79	0.14	106,106,107,107	0

6.5 Other polymers [i](#)

There are no such residues in this entry.