



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 12:02 pm GMT

PDB ID : 1RT5
Title : HIV-1 REVERSE TRANSCRIPTASE COMPLEXED WITH UC10
Authors : Ren, J.; Stammers, D.K.; Stuart, D.I.
Deposited on : 1998-07-29
Resolution : 2.90 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
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)	:	recalc28949

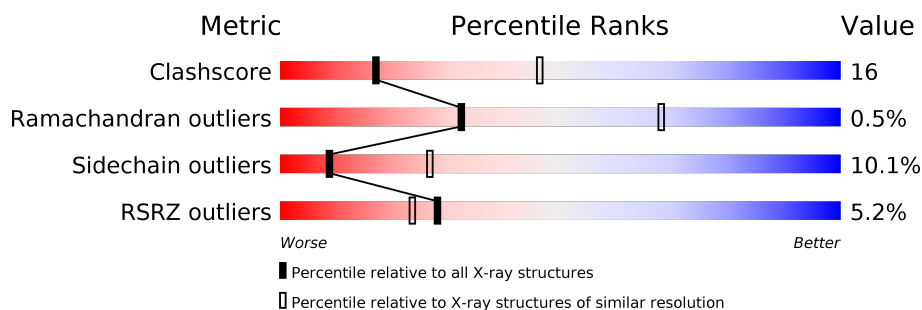
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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	112137	1807 (2.90-2.90)
Ramachandran outliers	110173	1768 (2.90-2.90)
Sidechain outliers	110143	1770 (2.90-2.90)
RSRZ outliers	101464	1596 (2.90-2.90)

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	560	
2	B	440	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	PO4	A	1001	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7780 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 HIV-1 REVERSE TRANSCRIPTASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	537	Total	C	N	O	S	0	0	0
			4396	2847	732	809	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	280	CSD	CYS	MODIFIED RESIDUE	UNP P04585

- Molecule 2 is a protein called HIV-1 REVERSE TRANSCRIPTASE.

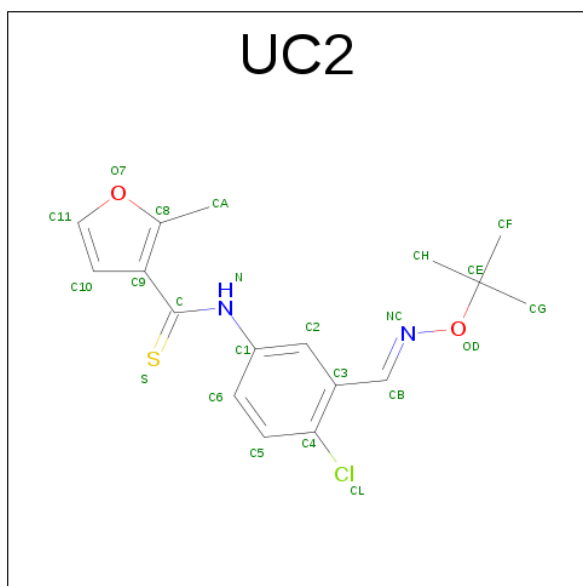
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	406	Total	C	N	O	S	0	0	0
			3351	2175	558	611	7			

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 4 is N-[4-CLORO-3-(T-BUTYLOXOME)PHENYL-2-METHYL-3-FURAN-CAR BOTHIAMIDE (three-letter code: UC2) (formula: C₁₇H₁₉ClN₂O₂S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	Cl	N	O	S	
			23	17	1	2	2	1	0
									0

- Molecule 5 is water.

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



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	140.30Å 111.00Å 73.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 2.90 29.65 – 2.90	Depositor EDS
% Data completeness (in resolution range)	86.1 (30.00-2.90) 87.2 (29.65-2.90)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.59 (at 2.90Å)	Xtriage
Refinement program	X-PLOR	Depositor
R, R_{free}	0.231 , 0.291 0.232 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	49.3	Xtriage
Anisotropy	0.163	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.30 , 83.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	7780	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.57% 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: CSD, PO4, UC2

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	0.34	0/4503	0.63	1/6117 (0.0%)
2	B	0.33	0/3442	0.62	0/4672
All	All	0.34	0/7945	0.63	1/10789 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	388	LYS	N-CA-C	-5.14	97.13	111.00

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4396	0	4451	151	0
2	B	3351	0	3386	93	0
3	A	5	0	0	0	0
4	A	23	0	19	5	0
5	A	4	0	0	1	0
5	B	1	0	0	0	0
All	All	7780	0	7856	243	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 (243) 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:34:LEU:HD11	1:A:62:ALA:HB2	1.27	1.16
1:A:358:ARG:HH11	1:A:358:ARG:HG2	1.16	1.07
1:A:540:LYS:HG2	2:B:280:CYS:SG	1.95	1.05
1:A:22:LYS:H	1:A:22:LYS:HD3	1.33	0.94
1:A:219:LYS:HD3	1:A:222:GLN:HE22	1.34	0.89
1:A:244:ILE:HG23	1:A:310:LEU:HD13	1.56	0.87
1:A:34:LEU:CD1	1:A:62:ALA:HB2	2.06	0.84
2:B:109:LEU:HG	2:B:216:THR:HG21	1.57	0.84
2:B:422:LEU:H	2:B:422:LEU:HD23	1.44	0.82
1:A:132:ILE:HB	1:A:142:ILE:HB	1.61	0.81
1:A:358:ARG:HG2	1:A:358:ARG:NH1	1.87	0.79
1:A:344:GLU:HB2	1:A:347:LYS:HD3	1.64	0.78
1:A:356:ARG:HH22	1:A:371:ALA:HB2	1.50	0.77
1:A:219:LYS:HD3	1:A:222:GLN:NE2	2.00	0.77
2:B:66:LYS:HG2	2:B:407:GLN:OE1	1.84	0.76
1:A:59:PRO:HG2	1:A:76:ASP:HB3	1.65	0.76
2:B:247:PRO:HB2	2:B:249:LYS:HZ1	1.49	0.76
1:A:193:LEU:HD13	1:A:197:GLN:HG3	1.68	0.76
2:B:242:GLN:HE21	2:B:242:GLN:HA	1.52	0.74
1:A:239:TRP:CZ2	1:A:316:GLY:HA3	2.22	0.74
2:B:84:THR:HG22	2:B:154:LYS:HE2	1.69	0.74
2:B:266:TRP:CD1	2:B:422:LEU:HD12	2.25	0.71
2:B:101:LYS:HD2	2:B:382:ILE:HG23	1.72	0.71
2:B:157:PRO:HG2	2:B:184:MET:HA	1.72	0.71
2:B:214:LEU:HD23	2:B:214:LEU:H	1.56	0.71
2:B:247:PRO:HB2	2:B:249:LYS:NZ	2.05	0.70
1:A:89:GLU:HB2	1:A:92:LEU:HD22	1.73	0.69
1:A:257:ILE:O	1:A:261:VAL:HG23	1.92	0.69
1:A:358:ARG:HH11	1:A:358:ARG:CG	2.02	0.69
2:B:114:ALA:HB2	2:B:214:LEU:HD13	1.75	0.68
1:A:235:HIS:HB2	1:A:238:LYS:O	1.93	0.67
2:B:354:TYR:HD1	2:B:374:LYS:HZ1	1.42	0.67
2:B:387:PRO:HG2	2:B:389:PHE:CE1	2.30	0.67
1:A:448:ARG:CZ	1:A:448:ARG:H	2.07	0.67
2:B:72:ARG:HH21	2:B:409:THR:HG22	1.60	0.66
1:A:340:GLN:HB3	1:A:351:THR:HG22	1.78	0.66
1:A:171:PHE:HB2	1:A:208:HIS:HD2	1.60	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:328:GLU:HG3	1:A:390:LYS:HB2	1.78	0.66
1:A:62:ALA:HB1	1:A:71:TRP:CE3	2.31	0.65
1:A:79:GLU:HG3	1:A:83:ARG:NH1	2.11	0.65
1:A:540:LYS:HB3	1:A:542:ILE:HD12	1.81	0.63
1:A:156:SER:HB2	1:A:157:PRO:HD3	1.81	0.62
2:B:109:LEU:CG	2:B:216:THR:HG21	2.29	0.61
2:B:158:ALA:O	2:B:161:GLN:HB2	2.01	0.61
1:A:171:PHE:HB2	1:A:208:HIS:CD2	2.35	0.61
1:A:22:LYS:H	1:A:22:LYS:CD	2.05	0.61
2:B:369:THR:HG22	2:B:398:TRP:CH2	2.36	0.60
2:B:263:LYS:HE2	2:B:425:LEU:HB3	1.84	0.60
2:B:109:LEU:HG	2:B:216:THR:CG2	2.31	0.60
2:B:57:ASN:HD22	2:B:143:ARG:NH1	2.00	0.60
1:A:62:ALA:HB1	1:A:71:TRP:HE3	1.66	0.59
1:A:448:ARG:HH11	1:A:448:ARG:HB3	1.67	0.59
1:A:17:ASP:O	1:A:83:ARG:HD3	2.03	0.59
1:A:31:ILE:O	1:A:35:VAL:HG23	2.03	0.59
1:A:542:ILE:HD12	1:A:542:ILE:H	1.66	0.59
4:A:999:UC2:C9	4:A:999:UC2:H2	2.34	0.58
1:A:344:GLU:CB	1:A:347:LYS:HD3	2.32	0.58
1:A:168:LEU:HD11	1:A:209:LEU:HD21	1.84	0.58
1:A:68:SER:C	1:A:70:LYS:H	2.06	0.57
2:B:206:ARG:HG3	2:B:216:THR:O	2.05	0.57
1:A:30:LYS:O	1:A:34:LEU:HD13	2.05	0.57
1:A:287:LYS:HG2	1:A:291:GLU:OE1	2.05	0.57
1:A:21:VAL:HG22	1:A:59:PRO:HD3	1.86	0.57
1:A:104:LYS:NZ	1:A:104:LYS:HB3	2.20	0.57
1:A:249:LYS:HG3	1:A:252:TRP:CE2	2.40	0.57
1:A:439:THR:HG21	2:B:289:LEU:HD13	1.86	0.56
2:B:120:LEU:HD12	2:B:150:PRO:HD3	1.87	0.56
1:A:50:ILE:HG13	1:A:143:ARG:HB3	1.88	0.56
1:A:216:THR:HG23	1:A:217:PRO:HD2	1.86	0.56
1:A:66:LYS:HD2	1:A:66:LYS:H	1.71	0.56
1:A:26:LEU:HD22	1:A:30:LYS:HE2	1.88	0.56
2:B:114:ALA:HB1	2:B:160:PHE:CZ	2.41	0.56
1:A:486:LEU:HB3	1:A:524:GLN:HG2	1.89	0.55
1:A:340:GLN:CB	1:A:351:THR:HG22	2.37	0.55
2:B:197:GLN:O	2:B:201:LYS:HB2	2.06	0.55
1:A:11:LYS:HB3	1:A:11:LYS:NZ	2.21	0.55
2:B:206:ARG:O	2:B:210:LEU:HD13	2.05	0.55
1:A:92:LEU:HD23	1:A:92:LEU:N	2.22	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:368:LEU:O	2:B:372:VAL:HG23	2.06	0.55
1:A:278:GLN:HG2	1:A:298:GLU:HB3	1.88	0.54
1:A:356:ARG:HB2	1:A:367:GLN:NE2	2.23	0.54
1:A:451:LYS:HG3	1:A:472:THR:O	2.08	0.54
1:A:432:GLU:HB2	1:A:433:PRO:HD2	1.91	0.53
1:A:122:GLU:CD	1:A:122:GLU:H	2.12	0.53
1:A:324:ASP:O	1:A:343:GLN:HG2	2.09	0.53
1:A:40:GLU:O	1:A:43:LYS:HG2	2.09	0.53
2:B:163:SER:O	2:B:167:ILE:HG13	2.08	0.53
1:A:366:LYS:O	1:A:369:THR:HB	2.09	0.53
1:A:13:LYS:HG2	1:A:83:ARG:O	2.09	0.53
2:B:319:TYR:HE1	2:B:325:LEU:HD13	1.74	0.53
2:B:395:LYS:O	2:B:399:GLU:HG2	2.08	0.52
1:A:129:ALA:HB1	1:A:143:ARG:NH1	2.24	0.52
2:B:325:LEU:HD21	2:B:383:TRP:CE3	2.45	0.52
1:A:296:THR:HG23	1:A:299:ALA:HB2	1.92	0.52
1:A:486:LEU:HB3	1:A:524:GLN:CG	2.40	0.52
1:A:516:GLU:HA	1:A:516:GLU:OE1	2.09	0.52
1:A:26:LEU:HD22	1:A:30:LYS:CE	2.41	0.51
1:A:540:LYS:CG	2:B:280:CYS:SG	2.86	0.51
2:B:360:ALA:HB2	2:B:366:LYS:HD2	1.90	0.51
1:A:37:ILE:O	1:A:40:GLU:HB3	2.10	0.51
1:A:94:ILE:HG23	1:A:229:TRP:CZ2	2.47	0.50
1:A:344:GLU:HG3	1:A:347:LYS:NZ	2.26	0.50
2:B:101:LYS:HG3	2:B:382:ILE:HA	1.94	0.50
1:A:125:ARG:HD3	1:A:147:ASN:HA	1.91	0.50
1:A:497:THR:O	1:A:535:TRP:HA	2.11	0.50
4:A:999:UC2:C10	4:A:999:UC2:C2	2.90	0.50
1:A:22:LYS:HD3	1:A:22:LYS:N	2.14	0.50
1:A:457:TYR:CE2	1:A:465:LYS:HB3	2.46	0.50
2:B:156:SER:HB2	2:B:157:PRO:HD3	1.93	0.50
2:B:23:GLN:OE1	2:B:59:PRO:HA	2.12	0.50
2:B:295:LEU:HB3	2:B:300:GLU:HG2	1.93	0.50
2:B:421:PRO:O	2:B:425:LEU:HD22	2.12	0.50
1:A:448:ARG:HB3	1:A:448:ARG:NH1	2.27	0.49
2:B:354:TYR:HD1	2:B:374:LYS:NZ	2.09	0.49
1:A:362:THR:HG22	1:A:363:ASN:H	1.78	0.49
1:A:252:TRP:CD1	1:A:295:LEU:HD12	2.48	0.49
4:A:999:UC2:HA1	4:A:999:UC2:S	2.53	0.49
1:A:454:LYS:HA	1:A:467:VAL:O	2.13	0.48
1:A:91:GLN:HG2	1:A:91:GLN:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:72:ARG:HH22	2:B:151:GLN:NE2	2.11	0.48
2:B:72:ARG:NH2	2:B:151:GLN:NE2	2.61	0.48
1:A:173:LYS:O	1:A:176:PRO:HD3	2.13	0.48
1:A:89:GLU:CB	1:A:92:LEU:HD22	2.42	0.48
2:B:125:ARG:NE	2:B:147:ASN:HA	2.27	0.48
2:B:109:LEU:CB	2:B:216:THR:HG21	2.44	0.48
2:B:288:ALA:O	2:B:291:GLU:HB3	2.14	0.48
2:B:365:VAL:O	2:B:369:THR:HG23	2.14	0.48
1:A:451:LYS:HD2	1:A:473:THR:HG22	1.95	0.48
1:A:405:TYR:CE2	1:A:407:GLN:HB3	2.49	0.48
1:A:94:ILE:HG23	1:A:229:TRP:HZ2	1.79	0.48
4:A:999:UC2:C9	4:A:999:UC2:C2	2.91	0.48
1:A:446:ALA:O	1:A:474:ASN:ND2	2.46	0.47
1:A:60:VAL:HG13	1:A:75:VAL:HG22	1.96	0.47
2:B:72:ARG:HH11	2:B:72:ARG:HG3	1.79	0.47
2:B:120:LEU:HD21	2:B:124:PHE:HD2	1.79	0.47
2:B:139:THR:HG22	2:B:140:PRO:HD2	1.95	0.47
1:A:126:LYS:HE3	1:A:127:TYR:CZ	2.50	0.47
2:B:115:TYR:OH	2:B:157:PRO:HG3	2.15	0.47
2:B:266:TRP:HZ3	2:B:426:TRP:CD1	2.32	0.47
1:A:13:LYS:HD3	1:A:84:THR:O	2.14	0.47
2:B:175:ASN:N	2:B:176:PRO:HD3	2.29	0.47
2:B:326:ILE:O	2:B:341:ILE:HA	2.14	0.47
2:B:66:LYS:O	2:B:67:ASP:HB2	2.15	0.47
1:A:122:GLU:HA	1:A:125:ARG:NE	2.31	0.46
1:A:344:GLU:HG3	1:A:347:LYS:HZ2	1.80	0.46
1:A:325:LEU:O	1:A:388:LYS:HG3	2.14	0.46
2:B:57:ASN:HD22	2:B:143:ARG:HH12	1.63	0.46
2:B:7:THR:HG22	2:B:119:PRO:HG2	1.97	0.46
1:A:56:TYR:O	1:A:143:ARG:NH1	2.44	0.46
1:A:68:SER:C	1:A:70:LYS:N	2.68	0.46
1:A:19:PRO:HG2	1:A:80:LEU:HB2	1.97	0.46
1:A:248:GLU:HA	1:A:307:ARG:NH2	2.30	0.46
1:A:448:ARG:HG2	1:A:449:GLU:H	1.81	0.46
1:A:406:TRP:CH2	2:B:418:ASN:HA	2.51	0.46
2:B:169:GLU:HB3	2:B:170:PRO:HD3	1.98	0.46
1:A:88:TRP:CE3	1:A:88:TRP:HA	2.51	0.46
2:B:261:VAL:HG13	2:B:276:VAL:HG21	1.98	0.46
1:A:356:ARG:HH22	1:A:371:ALA:CB	2.24	0.46
1:A:76:ASP:OD1	1:A:78:ARG:HG3	2.16	0.45
1:A:26:LEU:CA	1:A:30:LYS:HZ3	2.29	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:377:THR:O	1:A:381:VAL:HG23	2.16	0.45
1:A:446:ALA:O	1:A:448:ARG:NH2	2.50	0.45
1:A:13:LYS:H	1:A:13:LYS:HG2	1.53	0.45
2:B:115:TYR:HB3	2:B:149:LEU:HB2	1.99	0.45
2:B:63:ILE:H	2:B:63:ILE:HD13	1.82	0.45
1:A:393:ILE:HB	1:A:423:VAL:HG22	1.99	0.45
1:A:282:LEU:HD11	1:A:296:THR:HG22	1.98	0.45
2:B:28:GLU:HA	2:B:135:ILE:HD11	1.99	0.44
1:A:64:LYS:H	1:A:64:LYS:HG3	1.52	0.44
2:B:78:ARG:O	2:B:82:LYS:HG3	2.17	0.44
1:A:11:LYS:HB3	1:A:11:LYS:HZ3	1.82	0.44
2:B:426:TRP:O	2:B:429:LEU:HB2	2.18	0.44
1:A:116:PHE:O	1:A:148:VAL:HG21	2.17	0.44
1:A:300:GLU:HA	1:A:300:GLU:OE1	2.18	0.44
2:B:126:LYS:HG3	2:B:127:TYR:N	2.32	0.44
1:A:249:LYS:HG3	1:A:252:TRP:CZ2	2.52	0.44
1:A:282:LEU:HB3	1:A:293:ILE:HG21	1.98	0.44
2:B:270:ILE:O	2:B:272:PRO:HD3	2.18	0.44
1:A:361:HIS:ND1	1:A:361:HIS:N	2.65	0.43
1:A:448:ARG:HG2	1:A:449:GLU:N	2.33	0.43
1:A:72:ARG:HH21	1:A:74:LEU:CD1	2.32	0.43
2:B:422:LEU:H	2:B:422:LEU:CD2	2.18	0.43
1:A:26:LEU:HA	1:A:30:LYS:HZ3	1.82	0.43
1:A:542:ILE:H	1:A:542:ILE:CD1	2.26	0.43
1:A:184:MET:N	5:A:1004:HOH:O	2.51	0.43
1:A:226:PRO:HB3	1:A:235:HIS:CD2	2.53	0.43
1:A:277:ARG:HH11	1:A:277:ARG:HB2	1.84	0.43
2:B:328:GLU:O	2:B:339:TYR:HA	2.19	0.43
1:A:522:ILE:O	1:A:526:ILE:HG13	2.19	0.43
1:A:169:GLU:N	1:A:170:PRO:HD2	2.34	0.42
2:B:161:GLN:HA	2:B:161:GLN:NE2	2.33	0.42
2:B:359:GLY:HA2	2:B:361:HIS:CE1	2.54	0.42
2:B:94:ILE:HD12	2:B:94:ILE:N	2.34	0.42
1:A:179:VAL:O	1:A:189:VAL:HA	2.19	0.42
2:B:261:VAL:HG22	2:B:276:VAL:HG22	2.00	0.42
1:A:149:LEU:HA	1:A:150:PRO:HD3	1.80	0.42
1:A:296:THR:HG23	1:A:299:ALA:CB	2.49	0.42
1:A:358:ARG:H	1:A:358:ARG:HG3	1.49	0.42
1:A:109:LEU:HD23	1:A:216:THR:HG21	2.01	0.42
2:B:363:ASN:C	2:B:363:ASN:OD1	2.58	0.42
1:A:26:LEU:HB3	1:A:30:LYS:HB2	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:11:LYS:O	2:B:85:GLN:HG2	2.19	0.42
1:A:447:ASN:HA	1:A:448:ARG:NH2	2.35	0.42
2:B:287:LYS:HG2	2:B:287:LYS:H	1.69	0.42
2:B:104:LYS:HB2	2:B:192:ASP:HA	2.02	0.42
2:B:130:PHE:CZ	2:B:144:TYR:HB2	2.55	0.42
2:B:131:THR:OG1	2:B:143:ARG:HD2	2.19	0.42
2:B:87:PHE:CD1	2:B:87:PHE:N	2.88	0.42
1:A:184:MET:CE	1:A:184:MET:HA	2.50	0.42
1:A:463:ARG:NH1	1:A:488:ASP:O	2.52	0.41
1:A:469:LEU:HB2	1:A:472:THR:HG21	2.02	0.41
2:B:205:LEU:O	2:B:208:HIS:HB3	2.21	0.41
1:A:131:THR:OG1	1:A:143:ARG:HD3	2.20	0.41
2:B:94:ILE:HG22	2:B:94:ILE:O	2.20	0.41
1:A:181:TYR:CE2	1:A:183:TYR:HB2	2.54	0.41
1:A:234:LEU:HB3	4:A:999:UC2:CL	2.57	0.41
2:B:132:ILE:N	2:B:132:ILE:HD12	2.34	0.41
1:A:419:THR:O	1:A:419:THR:HG23	2.19	0.41
2:B:84:THR:CG2	2:B:154:LYS:HE2	2.46	0.41
1:A:103:LYS:HA	1:A:103:LYS:HD3	1.83	0.41
1:A:362:THR:HG23	1:A:366:LYS:HD3	2.01	0.41
2:B:30:LYS:HE2	2:B:30:LYS:HB3	1.88	0.41
1:A:115:TYR:O	1:A:149:LEU:HB2	2.21	0.41
1:A:225:PRO:HG3	1:A:227:PHE:CE2	2.55	0.41
2:B:183:TYR:HB3	2:B:188:TYR:HE2	1.85	0.41
1:A:447:ASN:HB3	1:A:450:THR:OG1	2.20	0.41
1:A:542:ILE:HG13	2:B:283:LEU:HB3	2.03	0.41
1:A:168:LEU:O	1:A:172:ARG:HG3	2.21	0.41
2:B:79:GLU:HA	2:B:82:LYS:HE2	2.03	0.41
1:A:116:PHE:C	1:A:148:VAL:HG21	2.41	0.41
1:A:135:ILE:HG13	1:A:135:ILE:H	1.58	0.41
1:A:244:ILE:HD11	1:A:263:LYS:HB3	2.03	0.41
1:A:401:TRP:CZ3	1:A:409:THR:HG21	2.56	0.41
2:B:81:ASN:OD1	2:B:153:TRP:HD1	2.05	0.40
1:A:21:VAL:CG2	1:A:59:PRO:HD3	2.51	0.40
1:A:96:HIS:CD2	1:A:98:ALA:H	2.40	0.40
2:B:206:ARG:HG2	2:B:210:LEU:CD1	2.51	0.40
2:B:345:PRO:O	2:B:346:PHE:HB2	2.21	0.40
1:A:271:TYR:O	1:A:272:PRO:C	2.59	0.40
1:A:328:GLU:O	1:A:339:TYR:HA	2.22	0.40
1:A:395:LYS:HD3	1:A:414:TRP:CZ2	2.57	0.40
2:B:353:LYS:HB2	2:B:353:LYS:HE3	1.93	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:66:LYS:HG3	2:B:66:LYS:H	1.60	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	532/560 (95%)	495 (93%)	34 (6%)	3 (1%)	28	64
2	B	400/440 (91%)	373 (93%)	25 (6%)	2 (0%)	32	68
All	All	932/1000 (93%)	868 (93%)	59 (6%)	5 (0%)	32	68

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	112	GLY
2	B	361	HIS
1	A	358	ARG
2	B	95	PRO
1	A	111	VAL

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.

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	481/499 (96%)	426 (89%)	55 (11%)	7	20
2	B	369/400 (92%)	338 (92%)	31 (8%)	13	36
All	All	850/899 (94%)	764 (90%)	86 (10%)	9	27

All (86) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	GLU
1	A	11	LYS
1	A	21	VAL
1	A	22	LYS
1	A	42	GLU
1	A	61	PHE
1	A	63	ILE
1	A	64	LYS
1	A	66	LYS
1	A	72	ARG
1	A	91	GLN
1	A	92	LEU
1	A	101	LYS
1	A	168	LEU
1	A	205	LEU
1	A	208	HIS
1	A	215	THR
1	A	218	ASP
1	A	222	GLN
1	A	230	MET
1	A	251	SER
1	A	279	LEU
1	A	283	LEU
1	A	289	LEU
1	A	290	THR
1	A	295	LEU
1	A	296	THR
1	A	297	GLU
1	A	301	LEU
1	A	303	LEU
1	A	311	LYS
1	A	312	GLU

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Mol	Chain	Res	Type
1	A	317	VAL
1	A	323	LYS
1	A	336	GLN
1	A	340	GLN
1	A	358	ARG
1	A	361	HIS
1	A	362	THR
1	A	368	LEU
1	A	424	LYS
1	A	448	ARG
1	A	451	LYS
1	A	452	LEU
1	A	471	ASP
1	A	475	GLN
1	A	480	GLN
1	A	491	LEU
1	A	493	VAL
1	A	496	VAL
1	A	513	SER
1	A	514	GLU
1	A	527	LYS
1	A	540	LYS
1	A	542	ILE
2	B	6	GLU
2	B	63	ILE
2	B	66	LYS
2	B	69	THR
2	B	87	PHE
2	B	101	LYS
2	B	109	LEU
2	B	123	ASP
2	B	139	THR
2	B	161	GLN
2	B	186	ASP
2	B	194	GLU
2	B	211	ARG
2	B	214	LEU
2	B	215	THR
2	B	242	GLN
2	B	283	LEU
2	B	287	LYS
2	B	303	LEU

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Mol	Chain	Res	Type
2	B	357	MET
2	B	358	ARG
2	B	361	HIS
2	B	362	THR
2	B	368	LEU
2	B	379	SER
2	B	399	GLU
2	B	418	ASN
2	B	422	LEU
2	B	424	LYS
2	B	425	LEU
2	B	432	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	174	GLN
1	A	222	GLN
1	A	235	HIS
1	A	242	GLN
1	A	315	HIS
1	A	332	GLN
1	A	373	GLN
2	B	57	ASN
2	B	85	GLN
2	B	151	GLN
2	B	161	GLN
2	B	242	GLN
2	B	269	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

1 non-standard protein/DNA/RNA residue is 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	$\# Z > 2$
1	CSD	A	280	1	4,7,8	1.47	1 (25%)	2,8,10	2.70	1 (50%)

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
1	CSD	A	280	1	-	1/2/6/8	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	280	CSD	CA-C	2.63	1.53	1.50

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	280	CSD	OD1-SG-CB	3.51	112.18	105.61

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	280	CSD	CA-CB-SG-OD1

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates

There are no carbohydrates in this entry.

5.6 Ligand geometry

2 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	PO4	A	1001	-	4,4,4	1.44	0	6,6,6	0.36	0
4	UC2	A	999	-	21,24,24	4.00	6 (28%)	21,34,34	1.55	3 (14%)

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
3	PO4	A	1001	-	-	0/0/0/0	0/0/0/0
4	UC2	A	999	-	-	1/10/15/15	0/1/2/2

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	999	UC2	OD-NC	-13.49	1.16	1.42
4	A	999	UC2	OD-CE	-4.67	1.40	1.47
4	A	999	UC2	C10-C11	-2.72	1.34	1.39
4	A	999	UC2	CA-C8	3.66	1.53	1.48
4	A	999	UC2	C-N	4.80	1.39	1.34
4	A	999	UC2	CB-NC	8.57	1.40	1.27

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	999	UC2	S-C-N	-4.63	117.91	125.72
4	A	999	UC2	C3-CB-NC	-3.37	116.33	121.16
4	A	999	UC2	C5-C4-C3	-2.35	119.85	121.94

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	999	UC2	S-C-N-C1

There are no ring outliers.

1 monomer is involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	999	UC2	5	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	536/560 (95%)	0.01	38 (7%) 17 12	11, 54, 124, 149	0
2	B	406/440 (92%)	-0.19	11 (2%) 55 50	10, 50, 108, 160	0
All	All	942/1000 (94%)	-0.08	49 (5%) 28 23	10, 52, 120, 160	0

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	68	SER	5.6
1	A	67	ASP	5.5
1	A	52	PRO	5.1
2	B	216	THR	5.0
1	A	30	LYS	4.5
1	A	29	GLU	4.3
2	B	6	GLU	4.1
1	A	28	GLU	4.1
1	A	132	ILE	4.0
1	A	450	THR	3.9
1	A	63	ILE	3.8
1	A	74	LEU	3.7
1	A	542	ILE	3.7
2	B	214	LEU	3.6
1	A	21	VAL	3.6
2	B	198	HIS	3.6
2	B	93	GLY	3.4
1	A	90	VAL	3.4
2	B	237	ASP	3.4
1	A	452	LEU	3.4
1	A	136	ASN	3.3
1	A	245	VAL	3.2
1	A	51	GLY	3.2
1	A	61	PHE	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	91	GLN	3.2
2	B	94	ILE	3.1
1	A	145	GLN	3.1
1	A	55	PRO	3.0
1	A	472	THR	2.7
2	B	197	GLN	2.6
1	A	1	PRO	2.6
2	B	67	ASP	2.6
1	A	72	ARG	2.6
1	A	130	PHE	2.5
1	A	133	PRO	2.5
1	A	64	LYS	2.5
2	B	215	THR	2.4
1	A	66	LYS	2.3
1	A	33	ALA	2.3
1	A	24	TRP	2.3
1	A	357	MET	2.2
1	A	134	SER	2.2
1	A	474	ASN	2.2
2	B	240	THR	2.1
1	A	2	ILE	2.1
1	A	92	LEU	2.1
1	A	541	GLY	2.0
1	A	448	ARG	2.0
1	A	32	LYS	2.0

6.2 Non-standard residues in protein, DNA, RNA chains [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(Å ²)	Q<0.9
1	CSD	A	280	8/9	0.95	0.12	-	40,44,51,60	0

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(Å ²)	Q<0.9
3	PO4	A	1001	5/5	0.93	0.27	3.39	100,101,108,109	0
4	UC2	A	999	23/23	0.98	0.12	-0.84	8,22,34,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.