



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:11 AM GMT

PDB ID : 3DOL
Title : Crystal structure of L100I mutant HIV-1 reverse transcriptase in complex with GW695634.
Authors : Chamberlain, P.P.; Ren, J.; Stammers, D.K.
Deposited on : 2008-07-04
Resolution : 2.50 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

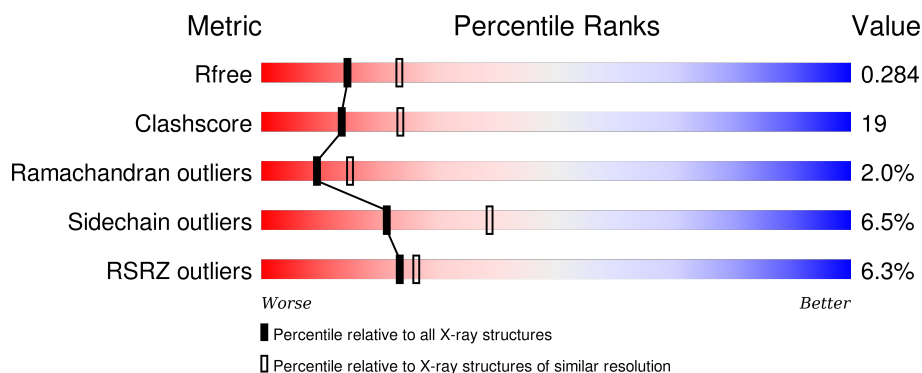
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	3553 (2.50-2.50)
Clashscore	102246	4242 (2.50-2.50)
Ramachandran outliers	100387	4156 (2.50-2.50)
Sidechain outliers	100360	4158 (2.50-2.50)
RSRZ outliers	91569	3562 (2.50-2.50)

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	<div> <div>6%</div> <div>60%32%</div> <div>• •</div> </div>
2	B	440	<div> <div>6%</div> <div>60%29%8%</div> <div>•</div> </div>

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	1300	-	-	-	X
3	PO4	A	1301	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 7902 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 Reverse transcriptase/ribonuclease H.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	537	Total	C	N	O	S	0	0	0
			4393	2843	731	811	8			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	100	ILE	LEU	ENGINEERED	UNP P04585

- Molecule 2 is a protein called p66 RT.

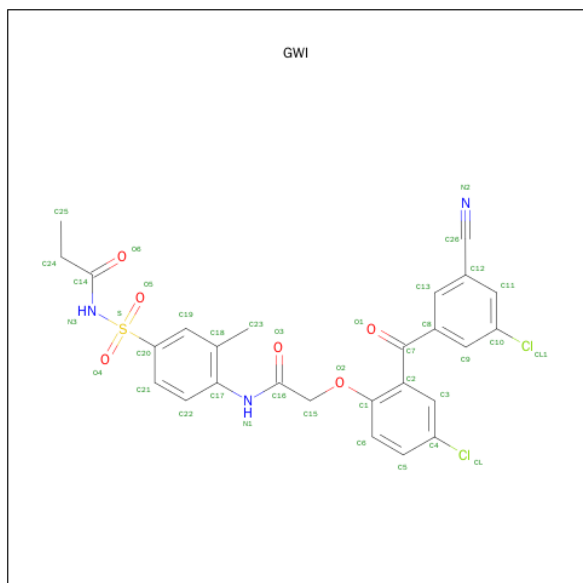
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	405	Total	C	N	O	S	0	0	0
			3346	2175	556	608	7			

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



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

- Molecule 4 is N-({4-[({4-CHLORO-2-[(3-CHLORO-5-CYANOPHENYL)CARBONYL]PHENOXY}ACETYL)AMINO]-3-METHYLPHENYL}SULFONYL)PROPANAMIDE (three-letter code: GWI) (formula: C₂₆H₂₁Cl₂N₃O₆S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C Cl N O S 38 26 2 3 6 1	0	0

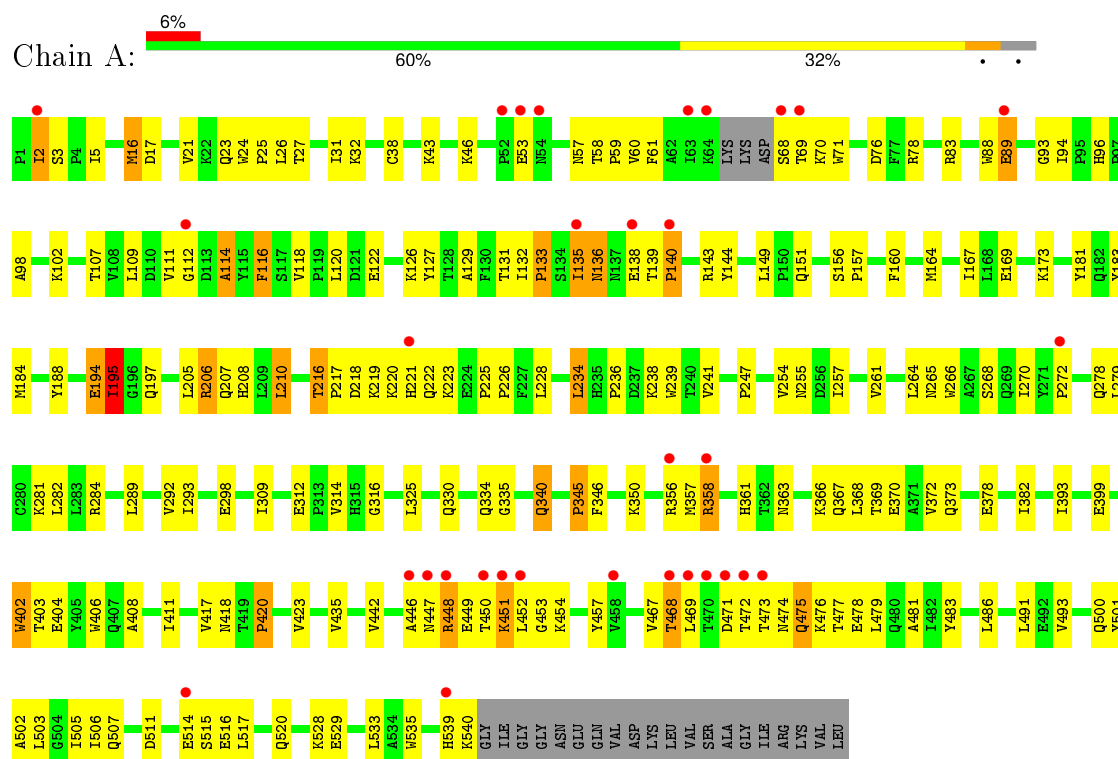
- Molecule 5 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	64	Total O 64 64	0	0
5	B	51	Total O 51 51	0	0

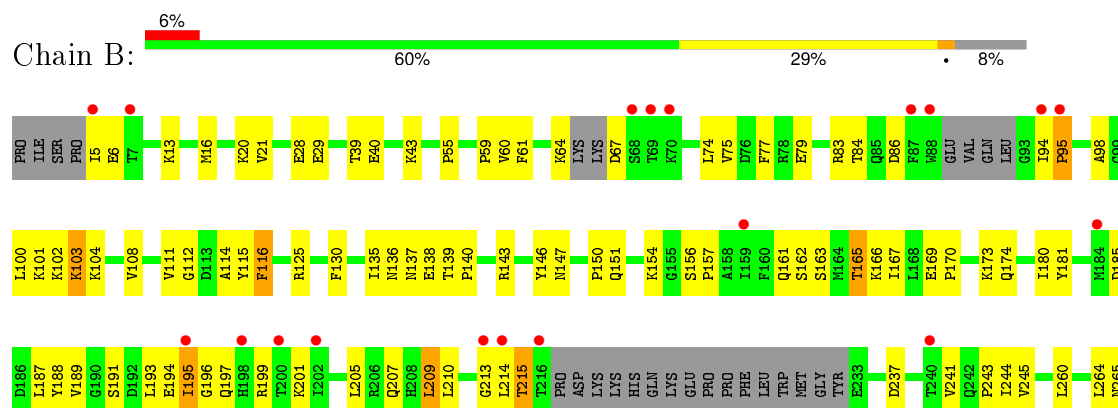
3 Residue-property plots

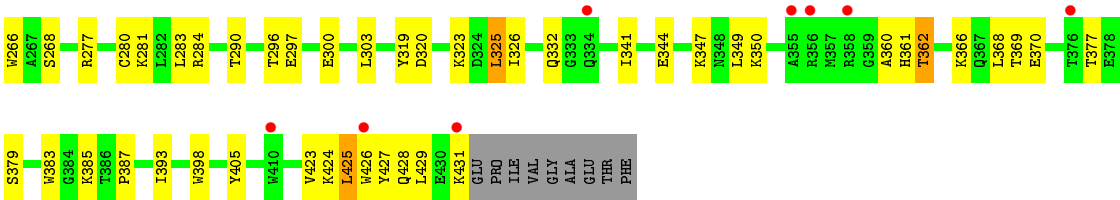
These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according 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 ($\text{RSRZ} > 2$). 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: Reverse transcriptase/ribonuclease H



• Molecule 2: p66 RT





4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	139.01Å 111.53Å 73.03Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.91 – 2.50 29.91 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.91-2.50) 100.0 (29.91-2.50)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.93 (at 2.51Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.224 , 0.288 0.222 , 0.284	Depositor DCC
R_{free} test set	1998 reflections (5.00%)	DCC
Wilson B-factor (Å ²)	60.7	Xtriage
Anisotropy	0.255	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 67.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 39995 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7902	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.12% 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: CSD, PO4, GWI

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	0.38	0/4501	0.65	1/6118 (0.0%)
2	B	0.39	0/3438	0.62	0/4669
All	All	0.39	0/7939	0.64	1/10787 (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	539	HIS	N-CA-C	6.32	128.06	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	4393	0	4435	182	0
2	B	3346	0	3374	118	0
3	A	10	0	0	0	0
4	A	38	0	21	0	0
5	A	64	0	0	9	0
5	B	51	0	0	2	0
All	All	7902	0	7830	290	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 (290) 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)
2:B:114:ALA:HB2	2:B:214:LEU:HD21	1.33	1.09
2:B:195:ILE:HD11	2:B:199:ARG:HE	1.30	0.96
1:A:358:ARG:HD3	1:A:358:ARG:H	1.30	0.95
1:A:96:HIS:HD2	1:A:98:ALA:H	1.16	0.92
2:B:61:PHE:CE1	2:B:74:LEU:HD23	2.04	0.91
1:A:135:ILE:H	1:A:135:ILE:HD12	1.35	0.90
1:A:218:ASP:O	1:A:222:GLN:HG3	1.71	0.90
2:B:5:ILE:HG22	2:B:6:GLU:H	1.40	0.87
2:B:244:ILE:HG13	2:B:426:TRP:CZ2	2.12	0.85
1:A:16:MET:HE2	1:A:83:ARG:HG2	1.58	0.84
1:A:217:PRO:HA	1:A:221:HIS:CE1	2.12	0.83
1:A:3:SER:HB3	1:A:5:ILE:HG13	1.59	0.82
2:B:114:ALA:HB2	2:B:214:LEU:CD2	2.09	0.81
1:A:206:ARG:HH11	1:A:216:THR:HG22	1.43	0.81
1:A:57:ASN:HD22	1:A:143:ARG:NH1	1.80	0.80
1:A:399:GLU:O	1:A:403:THR:HB	1.82	0.80
1:A:335:GLY:HA2	1:A:367:GLN:OE1	1.84	0.77
2:B:332:GLN:NE2	2:B:428:GLN:HB3	1.98	0.77
2:B:207:GLN:HA	2:B:210:LEU:HD12	1.68	0.74
2:B:215:THR:HG22	5:B:1115:HOH:O	1.86	0.74
1:A:451:LYS:HB3	1:A:471:ASP:HA	1.69	0.74
1:A:447:ASN:OD1	1:A:449:GLU:HB2	1.88	0.73
1:A:17:ASP:O	1:A:83:ARG:HD3	1.89	0.73
1:A:514:GLU:HG3	1:A:515:SER:N	2.04	0.73
2:B:320:ASP:HB3	2:B:323:LYS:HD2	1.71	0.72
2:B:169:GLU:HG3	2:B:173:LYS:NZ	2.05	0.72
2:B:197:GLN:O	2:B:201:LYS:HB2	1.90	0.72
1:A:46:LYS:HE2	1:A:116:PHE:HB3	1.73	0.71
2:B:5:ILE:HG22	2:B:6:GLU:N	2.06	0.70
1:A:57:ASN:HD22	1:A:143:ARG:HH12	1.38	0.69
2:B:161:GLN:O	2:B:165:THR:HG22	1.92	0.69
2:B:94:ILE:HG23	2:B:161:GLN:NE2	2.08	0.68
1:A:358:ARG:HD3	1:A:358:ARG:N	2.08	0.68
2:B:61:PHE:CZ	2:B:74:LEU:HD23	2.28	0.68
1:A:393:ILE:HB	1:A:423:VAL:HG22	1.76	0.67
1:A:69:THR:HG22	1:A:69:THR:O	1.95	0.67
2:B:125:ARG:HE	2:B:147:ASN:HA	1.59	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:402:TRP:CH2	2:B:362:THR:HA	2.30	0.67
1:A:96:HIS:CD2	1:A:98:ALA:H	2.05	0.67
2:B:344:GLU:HB2	2:B:347:LYS:HD2	1.77	0.66
2:B:180:ILE:HG12	2:B:189:VAL:HG12	1.76	0.65
2:B:195:ILE:HD11	2:B:199:ARG:NE	2.07	0.65
1:A:206:ARG:NH1	1:A:216:THR:HG22	2.09	0.65
1:A:216:THR:HG23	1:A:217:PRO:N	2.12	0.65
2:B:332:GLN:HE22	2:B:428:GLN:HB3	1.63	0.63
1:A:447:ASN:HB3	1:A:450:THR:OG1	1.98	0.63
1:A:151:GLN:HB3	5:A:1007:HOH:O	1.97	0.63
1:A:435:VAL:HG22	2:B:290:THR:HG21	1.81	0.63
1:A:241:VAL:HG21	1:A:270:ILE:HD13	1.81	0.62
1:A:122:GLU:H	1:A:122:GLU:CD	2.01	0.62
2:B:100:LEU:HD22	2:B:181:TYR:HB3	1.82	0.62
2:B:205:LEU:O	2:B:209:LEU:HD22	2.00	0.62
1:A:292:VAL:C	1:A:293:ILE:HD12	2.20	0.62
1:A:206:ARG:HG2	1:A:206:ARG:HH11	1.65	0.62
2:B:243:PRO:HG2	2:B:431:LYS:HB3	1.82	0.62
1:A:448:ARG:H	1:A:448:ARG:HD3	1.64	0.62
1:A:219:LYS:HD2	1:A:219:LYS:H	1.64	0.61
1:A:27:THR:O	1:A:31:ILE:HG13	2.02	0.60
2:B:281:LYS:O	2:B:284:ARG:HG3	2.02	0.60
2:B:98:ALA:HA	2:B:101:LYS:HE3	1.84	0.59
2:B:209:LEU:HD12	2:B:214:LEU:CD1	2.33	0.59
1:A:358:ARG:CD	1:A:358:ARG:H	2.10	0.59
2:B:245:VAL:HG13	2:B:431:LYS:HB2	1.84	0.59
1:A:393:ILE:HB	1:A:423:VAL:CG2	2.31	0.59
1:A:514:GLU:HG3	1:A:515:SER:H	1.67	0.59
2:B:170:PRO:O	2:B:174:GLN:HG3	2.02	0.59
2:B:209:LEU:HB3	2:B:214:LEU:HB2	1.84	0.59
1:A:118:VAL:HB	1:A:149:LEU:HD22	1.85	0.58
1:A:239:TRP:CE2	1:A:316:GLY:HA3	2.39	0.58
2:B:169:GLU:HB3	2:B:170:PRO:HD3	1.84	0.58
2:B:281:LYS:HG2	2:B:284:ARG:HH21	1.69	0.58
1:A:102:LYS:HE2	1:A:236:PRO:O	2.04	0.57
2:B:195:ILE:HG23	2:B:196:GLY:H	1.69	0.57
2:B:163:SER:O	2:B:167:ILE:HG23	2.05	0.57
1:A:257:ILE:O	1:A:261:VAL:HG23	2.05	0.57
1:A:402:TRP:CZ2	2:B:362:THR:HA	2.40	0.57
1:A:452:LEU:HA	1:A:469:LEU:O	2.05	0.57
2:B:195:ILE:HG23	2:B:196:GLY:N	2.19	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:475:GLN:HG3	1:A:476:LYS:N	2.19	0.56
1:A:451:LYS:O	1:A:471:ASP:N	2.39	0.56
1:A:218:ASP:OD1	1:A:219:LYS:N	2.39	0.56
2:B:320:ASP:CB	2:B:323:LYS:HD2	2.36	0.56
1:A:23:GLN:HG3	1:A:133:PRO:HG3	1.85	0.56
2:B:319:TYR:OH	2:B:385:LYS:HD3	2.06	0.56
1:A:446:ALA:CB	1:A:453:GLY:HA3	2.36	0.56
1:A:78:ARG:HD2	5:A:1111:HOH:O	2.05	0.56
1:A:23:GLN:CG	1:A:133:PRO:HG3	2.36	0.56
1:A:363:ASN:HA	1:A:511:ASP:OD2	2.06	0.56
1:A:293:ILE:HD12	1:A:293:ILE:N	2.21	0.55
1:A:136:ASN:O	1:A:139:THR:HG22	2.06	0.55
2:B:244:ILE:HG13	2:B:426:TRP:CH2	2.42	0.55
2:B:111:VAL:HG22	2:B:185:ASP:O	2.05	0.54
1:A:446:ALA:HB1	1:A:452:LEU:O	2.07	0.54
2:B:266:TRP:HZ3	2:B:426:TRP:CE3	2.25	0.54
2:B:169:GLU:HG3	2:B:173:LYS:HZ2	1.72	0.54
1:A:38:CYS:SG	1:A:132:ILE:HD11	2.47	0.54
2:B:21:VAL:HB	2:B:59:PRO:HD3	1.90	0.54
1:A:156:SER:HB2	1:A:157:PRO:HD3	1.88	0.54
2:B:13:LYS:HB2	2:B:16:MET:HG3	1.90	0.54
1:A:417:VAL:O	1:A:417:VAL:HG13	2.07	0.54
2:B:166:LYS:NZ	2:B:166:LYS:HB2	2.23	0.54
2:B:125:ARG:HG2	2:B:146:TYR:O	2.08	0.53
1:A:408:ALA:HB3	2:B:393:ILE:HG13	1.90	0.53
1:A:270:ILE:CG2	1:A:314:VAL:HG21	2.38	0.53
2:B:241:VAL:HG22	2:B:350:LYS:HA	1.91	0.53
1:A:491:LEU:HB3	1:A:529:GLU:HG3	1.91	0.53
2:B:213:GLY:O	2:B:214:LEU:HG	2.09	0.53
1:A:46:LYS:NZ	1:A:46:LYS:HB3	2.24	0.53
1:A:257:ILE:HD13	1:A:282:LEU:HD23	1.91	0.52
1:A:194:GLU:OE1	1:A:195:ILE:HG23	2.10	0.52
1:A:194:GLU:OE2	1:A:195:ILE:HG22	2.09	0.52
1:A:448:ARG:CD	1:A:448:ARG:H	2.23	0.52
1:A:442:VAL:HG12	1:A:457:TYR:HB3	1.92	0.52
1:A:234:LEU:HD22	1:A:234:LEU:N	2.25	0.52
1:A:228:LEU:HD22	1:A:228:LEU:N	2.26	0.51
1:A:366:LYS:O	1:A:369:THR:HB	2.09	0.51
1:A:94:ILE:HD12	1:A:183:TYR:HE2	1.73	0.51
1:A:221:HIS:ND1	1:A:221:HIS:O	2.44	0.51
1:A:514:GLU:CG	1:A:515:SER:N	2.73	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:457:TYR:C	1:A:457:TYR:CD1	2.83	0.51
2:B:325:LEU:HD21	2:B:383:TRP:CE3	2.45	0.51
1:A:502:ALA:HA	1:A:505:ILE:HD12	1.92	0.51
1:A:2:ILE:CG2	1:A:3:SER:N	2.73	0.51
1:A:96:HIS:HE1	5:A:1015:HOH:O	1.93	0.51
2:B:426:TRP:O	2:B:429:LEU:HB2	2.10	0.51
2:B:366:LYS:O	2:B:370:GLU:HG3	2.10	0.51
2:B:266:TRP:CZ2	2:B:423:VAL:HG22	2.46	0.51
1:A:83:ARG:HG3	1:A:83:ARG:HH11	1.75	0.51
1:A:3:SER:CB	1:A:5:ILE:HG13	2.37	0.51
2:B:28:GLU:HG3	2:B:135:ILE:HD11	1.91	0.51
1:A:23:GLN:OE1	1:A:60:VAL:HG12	2.11	0.51
1:A:279:LEU:O	1:A:282:LEU:HB2	2.11	0.50
1:A:418:ASN:O	1:A:420:PRO:HD3	2.11	0.50
1:A:451:LYS:O	1:A:472:THR:N	2.41	0.50
2:B:103:LYS:HE3	2:B:191:SER:HA	1.93	0.50
2:B:425:LEU:HA	2:B:428:GLN:HE21	1.76	0.49
1:A:483:TYR:CE1	1:A:520:GLN:HB3	2.47	0.49
1:A:350:LYS:HE3	1:A:378:GLU:OE2	2.10	0.49
2:B:379:SER:CB	2:B:387:PRO:HD3	2.42	0.49
1:A:478:GLU:O	1:A:481:ALA:HB3	2.12	0.49
1:A:418:ASN:C	1:A:420:PRO:HD3	2.33	0.49
2:B:125:ARG:NE	2:B:147:ASN:HA	2.25	0.49
1:A:225:PRO:HA	1:A:226:PRO:C	2.33	0.49
1:A:454:LYS:HE2	5:A:1085:HOH:O	2.13	0.49
1:A:314:VAL:HG23	1:A:314:VAL:O	2.13	0.49
1:A:206:ARG:NH1	1:A:206:ARG:HG2	2.27	0.49
1:A:334:GLN:HB3	1:A:356:ARG:HH21	1.78	0.49
1:A:270:ILE:HG21	1:A:314:VAL:HG21	1.95	0.49
1:A:473:THR:OG1	1:A:476:LYS:HG3	2.13	0.49
2:B:210:LEU:HD21	2:B:215:THR:O	2.14	0.48
2:B:108:VAL:HG22	2:B:188:TYR:CE2	2.48	0.48
2:B:426:TRP:CZ3	2:B:427:TYR:CE1	3.01	0.48
1:A:129:ALA:HA	1:A:144:TYR:O	2.13	0.48
1:A:23:GLN:O	1:A:25:PRO:HD3	2.14	0.48
2:B:210:LEU:CD2	2:B:215:THR:O	2.62	0.48
1:A:57:ASN:HA	1:A:129:ALA:O	2.13	0.48
1:A:506:ILE:HD11	1:A:533:LEU:HG	1.95	0.48
1:A:222:GLN:HB3	1:A:223:LYS:H	1.30	0.47
1:A:38:CYS:HB3	1:A:144:TYR:CE1	2.49	0.47
2:B:245:VAL:CG1	2:B:431:LYS:HB2	2.43	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:68:SER:OG	1:A:70:LYS:HG2	2.14	0.47
2:B:5:ILE:CG2	2:B:6:GLU:H	2.09	0.47
1:A:24:TRP:CE3	1:A:61:PHE:HZ	2.32	0.47
1:A:402:TRP:CD1	1:A:402:TRP:C	2.88	0.47
1:A:451:LYS:NZ	1:A:471:ASP:OD2	2.45	0.47
2:B:349:LEU:O	2:B:350:LYS:HD2	2.15	0.47
2:B:369:THR:HG22	2:B:398:TRP:CH2	2.48	0.47
5:A:1103:HOH:O	2:B:280:CYS:SG	2.61	0.47
1:A:16:MET:CE	1:A:83:ARG:HG2	2.39	0.47
1:A:58:THR:CG2	1:A:59:PRO:HD2	2.44	0.47
1:A:278:GLN:O	1:A:282:LEU:HD13	2.13	0.47
2:B:112:GLY:HA3	2:B:151:GLN:OE1	2.13	0.47
1:A:181:TYR:HB2	1:A:188:TYR:HB2	1.95	0.47
1:A:475:GLN:O	1:A:479:LEU:HG	2.15	0.47
2:B:84:THR:HB	2:B:154:LYS:HE2	1.97	0.47
2:B:115:TYR:OH	2:B:157:PRO:HB3	2.14	0.47
2:B:296:THR:O	2:B:300:GLU:HG2	2.15	0.47
1:A:216:THR:CG2	1:A:217:PRO:N	2.78	0.47
1:A:57:ASN:ND2	1:A:143:ARG:NH1	2.58	0.47
1:A:136:ASN:N	1:A:136:ASN:OD1	2.48	0.47
1:A:76:ASP:OD2	1:A:78:ARG:HG3	2.16	0.46
1:A:126:LYS:NZ	1:A:127:TYR:CZ	2.83	0.46
2:B:146:TYR:CD2	2:B:150:PRO:HB3	2.51	0.46
2:B:360:ALA:O	2:B:362:THR:N	2.47	0.46
1:A:278:GLN:HA	1:A:281:LYS:HG3	1.98	0.46
1:A:58:THR:HG23	1:A:59:PRO:HD2	1.97	0.46
2:B:98:ALA:HA	2:B:101:LYS:CE	2.44	0.46
1:A:135:ILE:H	1:A:135:ILE:CD1	2.09	0.46
1:A:216:THR:HB	5:A:1043:HOH:O	2.15	0.46
1:A:3:SER:HB3	1:A:5:ILE:CG1	2.40	0.46
2:B:277:ARG:HG2	2:B:277:ARG:HH21	1.79	0.46
2:B:297:GLU:O	2:B:300:GLU:HB2	2.15	0.46
1:A:265:ASN:O	1:A:268:SER:HB2	2.15	0.46
1:A:357:MET:O	1:A:367:GLN:NE2	2.49	0.46
2:B:209:LEU:HD12	2:B:214:LEU:HD13	1.97	0.46
1:A:403:THR:CG2	1:A:404:GLU:N	2.78	0.46
1:A:516:GLU:HG3	1:A:520:GLN:HE21	1.81	0.46
2:B:194:GLU:O	2:B:195:ILE:C	2.55	0.46
2:B:98:ALA:HA	2:B:101:LYS:NZ	2.30	0.46
2:B:325:LEU:HD12	2:B:325:LEU:HA	1.84	0.46
1:A:446:ALA:HB2	1:A:477:THR:HG21	1.97	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:GLU:HG2	5:A:1011:HOH:O	2.15	0.45
1:A:131:THR:OG1	1:A:143:ARG:HD2	2.16	0.45
1:A:403:THR:HG23	1:A:404:GLU:OE1	2.15	0.45
2:B:100:LEU:HD22	2:B:181:TYR:CB	2.46	0.45
1:A:114:ALA:HB1	1:A:160:PHE:CE2	2.50	0.45
1:A:89:GLU:OE1	1:A:89:GLU:O	2.35	0.45
1:A:473:THR:HB	5:A:1086:HOH:O	2.16	0.45
2:B:426:TRP:HZ3	2:B:427:TYR:CE1	2.33	0.45
2:B:98:ALA:O	2:B:102:LYS:HE2	2.16	0.45
2:B:64:LYS:C	2:B:67:ASP:N	2.69	0.45
1:A:403:THR:HG23	1:A:404:GLU:N	2.31	0.45
1:A:514:GLU:CG	1:A:515:SER:H	2.28	0.45
2:B:103:LYS:HB3	2:B:104:LYS:H	1.67	0.45
1:A:88:TRP:CE2	2:B:143:ARG:HD3	2.52	0.45
1:A:221:HIS:CG	1:A:221:HIS:O	2.69	0.45
2:B:139:THR:HG22	2:B:140:PRO:O	2.16	0.45
1:A:46:LYS:HB3	1:A:46:LYS:HZ3	1.82	0.44
2:B:156:SER:N	2:B:157:PRO:HD2	2.32	0.44
1:A:417:VAL:O	1:A:417:VAL:CG1	2.64	0.44
2:B:108:VAL:HG22	2:B:188:TYR:CD2	2.51	0.44
1:A:2:ILE:HG23	1:A:3:SER:N	2.32	0.44
2:B:424:LYS:C	2:B:424:LYS:HD3	2.38	0.44
1:A:88:TRP:CZ2	2:B:143:ARG:HD3	2.52	0.44
2:B:139:THR:HG23	2:B:140:PRO:HD2	1.98	0.44
2:B:39:THR:O	2:B:43:LYS:HG2	2.18	0.44
1:A:216:THR:HG23	1:A:217:PRO:CD	2.48	0.44
1:A:167:ILE:O	1:A:208:HIS:HE1	2.00	0.44
2:B:194:GLU:OE1	2:B:195:ILE:HG22	2.18	0.43
1:A:454:LYS:HB3	1:A:468:THR:HA	2.00	0.43
1:A:503:LEU:HA	1:A:506:ILE:HD12	1.99	0.43
2:B:360:ALA:C	2:B:362:THR:H	2.22	0.43
2:B:61:PHE:HE1	2:B:74:LEU:HD23	1.72	0.43
2:B:94:ILE:HB	2:B:95:PRO:HD2	1.99	0.43
1:A:257:ILE:CD1	1:A:282:LEU:HD23	2.48	0.43
2:B:75:VAL:HG11	2:B:77:PHE:CZ	2.53	0.43
1:A:281:LYS:O	1:A:284:ARG:HG3	2.18	0.43
1:A:111:VAL:HG21	1:A:164:MET:HE2	1.99	0.43
1:A:330:GLN:NE2	1:A:340:GLN:OE1	2.51	0.43
2:B:116:PHE:C	2:B:116:PHE:CD1	2.91	0.43
2:B:79:GLU:O	2:B:83:ARG:HG3	2.17	0.43
2:B:111:VAL:HB	2:B:214:LEU:HD22	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:GLY:O	1:A:356:ARG:HA	2.19	0.43
1:A:402:TRP:CH2	2:B:362:THR:HG22	2.54	0.43
1:A:132:ILE:O	1:A:133:PRO:O	2.36	0.42
1:A:183:TYR:CE1	1:A:184:MET:HG3	2.53	0.42
1:A:210:LEU:O	1:A:210:LEU:HD13	2.19	0.42
1:A:370:GLU:O	1:A:373:GLN:HG3	2.19	0.42
2:B:111:VAL:HG11	2:B:187:LEU:HD22	2.01	0.42
1:A:206:ARG:CG	1:A:216:THR:HG21	2.49	0.42
1:A:393:ILE:HD12	1:A:423:VAL:HG23	2.00	0.42
1:A:69:THR:CG2	1:A:69:THR:O	2.66	0.42
1:A:503:LEU:HD13	1:A:535:TRP:CB	2.49	0.42
1:A:278:GLN:HG3	1:A:298:GLU:HB3	2.02	0.42
1:A:408:ALA:HB3	2:B:393:ILE:CG1	2.48	0.42
1:A:454:LYS:HD3	1:A:468:THR:HG22	2.01	0.42
2:B:344:GLU:CB	2:B:347:LYS:HD2	2.48	0.42
1:A:107:THR:HG22	1:A:109:LEU:HD12	2.02	0.42
2:B:162:SER:O	2:B:165:THR:HG23	2.20	0.42
2:B:277:ARG:O	2:B:281:LYS:HG3	2.20	0.42
1:A:93:GLY:HA3	2:B:137:ASN:ND2	2.35	0.42
2:B:326:ILE:O	2:B:341:ILE:HA	2.19	0.42
1:A:219:LYS:HD2	1:A:219:LYS:N	2.33	0.42
1:A:293:ILE:CD1	1:A:293:ILE:N	2.83	0.42
2:B:320:ASP:C	2:B:320:ASP:OD1	2.59	0.41
1:A:46:LYS:CB	1:A:46:LYS:NZ	2.84	0.41
1:A:24:TRP:CE3	1:A:61:PHE:CZ	3.08	0.41
1:A:181:TYR:CD2	2:B:138:GLU:HG3	2.55	0.41
1:A:210:LEU:HD23	5:A:1041:HOH:O	2.19	0.41
1:A:372:VAL:HG11	1:A:411:ILE:HG23	2.02	0.41
1:A:475:GLN:NE2	1:A:501:TYR:CE2	2.89	0.41
1:A:378:GLU:O	1:A:382:ILE:HG12	2.21	0.41
1:A:454:LYS:CB	1:A:467:VAL:O	2.69	0.41
2:B:60:VAL:HG11	2:B:130:PHE:CD1	2.56	0.41
1:A:345:PRO:O	1:A:346:PHE:HB2	2.19	0.41
2:B:265:ASN:O	2:B:268:SER:OG	2.39	0.41
1:A:5:ILE:H	1:A:5:ILE:HG13	1.64	0.41
1:A:451:LYS:CG	1:A:472:THR:O	2.68	0.41
1:A:94:ILE:HD12	1:A:183:TYR:CE2	2.55	0.41
1:A:32:LYS:HA	1:A:32:LYS:HD3	1.91	0.41
1:A:309:ILE:O	1:A:312:GLU:HB2	2.21	0.41
1:A:70:LYS:HG3	1:A:71:TRP:N	2.35	0.40
1:A:406:TRP:CD1	1:A:507:GLN:HG2	2.57	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:254:VAL:HB	1:A:289:LEU:O	2.21	0.40
1:A:486:LEU:O	1:A:528:LYS:NZ	2.52	0.40
2:B:377:THR:HG23	5:B:1100:HOH:O	2.22	0.40
1:A:218:ASP:O	1:A:222:GLN:CG	2.56	0.40
2:B:136:ASN:O	2:B:137:ASN:HB2	2.22	0.40
1:A:107:THR:HG22	1:A:109:LEU:CD1	2.52	0.40
2:B:260:LEU:O	2:B:264:LEU:HG	2.20	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%)	481 (90%)	38 (7%)	13 (2%)	7	11
2	B	397/440 (90%)	362 (91%)	29 (7%)	6 (2%)	13	22
All	All	929/1000 (93%)	843 (91%)	67 (7%)	19 (2%)	9	15

All (19) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	114	ALA
1	A	133	PRO
2	B	103	LYS
2	B	193	LEU
2	B	361	HIS
1	A	53	GLU
1	A	112	GLY
1	A	140	PRO
1	A	272	PRO
2	B	195	ILE
1	A	247	PRO

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Mol	Chain	Res	Type
1	A	116	PHE
1	A	138	GLU
2	B	237	ASP
1	A	345	PRO
1	A	195	ILE
1	A	361	HIS
2	B	95	PRO
1	A	420	PRO

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	481/499 (96%)	442 (92%)	39 (8%)	15	27
2	B	368/400 (92%)	352 (96%)	16 (4%)	35	61
All	All	849/899 (94%)	794 (94%)	55 (6%)	21	39

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	ILE
1	A	16	MET
1	A	21	VAL
1	A	26	LEU
1	A	43	LYS
1	A	89	GLU
1	A	120	LEU
1	A	135	ILE
1	A	136	ASN
1	A	140	PRO
1	A	173	LYS
1	A	194	GLU
1	A	195	ILE
1	A	197	GLN
1	A	205	LEU

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Mol	Chain	Res	Type
1	A	206	ARG
1	A	207	GLN
1	A	210	LEU
1	A	216	THR
1	A	220	LYS
1	A	234	LEU
1	A	238	LYS
1	A	255	ASN
1	A	264	LEU
1	A	266	TRP
1	A	325	LEU
1	A	340	GLN
1	A	358	ARG
1	A	368	LEU
1	A	402	TRP
1	A	448	ARG
1	A	451	LYS
1	A	468	THR
1	A	474	ASN
1	A	475	GLN
1	A	493	VAL
1	A	500	GLN
1	A	517	LEU
1	A	540	LYS
2	B	20	LYS
2	B	29	GLU
2	B	40	GLU
2	B	55	PRO
2	B	86	ASP
2	B	116	PHE
2	B	165	THR
2	B	209	LEU
2	B	215	THR
2	B	283	LEU
2	B	303	LEU
2	B	325	LEU
2	B	362	THR
2	B	368	LEU
2	B	405	TYR
2	B	425	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	57	ASN
1	A	96	HIS
1	A	147	ASN
1	A	174	GLN
1	A	207	GLN
1	A	208	HIS
1	A	221	HIS
1	A	222	GLN
1	A	242	GLN
1	A	255	ASN
1	A	278	GLN
1	A	336	GLN
1	A	474	ASN
1	A	475	GLN
1	A	487	GLN
1	A	500	GLN
1	A	520	GLN
1	A	524	GLN
2	B	57	ASN
2	B	85	GLN
2	B	137	ASN
2	B	147	ASN
2	B	161	GLN
2	B	174	GLN
2	B	182	GLN
2	B	197	GLN
2	B	207	GLN
2	B	208	HIS
2	B	235	HIS
2	B	278	GLN
2	B	332	GLN
2	B	428	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	3,7,8	0.94	0	3,8,10	4.61	1 (33%)

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

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	A	280	CSD	OD1-SG-CB	7.76	118.33	105.40

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 [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

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	PO4	A	1300	-	4,4,4	1.14	0	6,6,6	0.27	0
3	PO4	A	1301	-	4,4,4	1.16	0	6,6,6	0.27	0
4	GWI	A	999	-	39,40,40	2.68	13 (33%)	54,57,57	1.52	10 (18%)

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	1300	-	-	0/0/0/0	0/0/0/0
3	PO4	A	1301	-	-	0/0/0/0	0/0/0/0
4	GWI	A	999	-	-	0/32/32/32	0/3/3/3

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	999	GWI	C17-N1	-2.07	1.37	1.41
4	A	999	GWI	C11-C12	2.01	1.43	1.39
4	A	999	GWI	C8-C7	2.15	1.53	1.49
4	A	999	GWI	C11-C10	2.29	1.42	1.38
4	A	999	GWI	C9-C8	2.31	1.42	1.39
4	A	999	GWI	C13-C8	2.53	1.43	1.39
4	A	999	GWI	O4-S	3.56	1.47	1.43
4	A	999	GWI	C22-C21	5.45	1.48	1.38
4	A	999	GWI	C3-C4	5.51	1.47	1.38
4	A	999	GWI	C2-C1	5.52	1.51	1.40
4	A	999	GWI	C19-C20	5.92	1.50	1.39
4	A	999	GWI	C5-C6	6.14	1.49	1.38
4	A	999	GWI	C17-C18	6.82	1.53	1.40

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	A	999	GWI	O5-S-O4	-5.89	111.72	119.54
4	A	999	GWI	C22-C17-C18	-3.04	117.20	120.67
4	A	999	GWI	C20-S-N3	-2.44	102.77	105.93
4	A	999	GWI	O2-C15-C16	-2.21	104.41	110.83
4	A	999	GWI	C1-C2-C7	-2.05	119.34	122.78
4	A	999	GWI	C15-C16-N1	2.01	119.36	115.77
4	A	999	GWI	C2-C7-C8	2.04	122.94	119.47
4	A	999	GWI	C8-C9-C10	2.10	121.52	119.19
4	A	999	GWI	C15-O2-C1	2.44	123.16	117.58
4	A	999	GWI	O5-S-C20	2.95	111.69	107.96

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	536/560 (95%)	0.15	32 (5%) 25 28	30, 61, 108, 147	0
2	B	405/440 (92%)	0.19	27 (6%) 21 23	32, 60, 109, 154	0
All	All	941/1000 (94%)	0.16	59 (6%) 23 26	30, 61, 109, 154	0

All (59) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	452	LEU	8.9
1	A	446	ALA	7.6
1	A	470	THR	6.5
2	B	88	TRP	6.5
1	A	469	LEU	5.9
1	A	539	HIS	5.1
2	B	240	THR	4.6
2	B	358	ARG	4.5
2	B	198	HIS	4.5
1	A	64	LYS	4.4
1	A	450	THR	4.2
2	B	69	THR	4.1
1	A	140	PRO	4.1
1	A	68	SER	4.0
2	B	213	GLY	4.0
2	B	356	ARG	3.9
1	A	52	PRO	3.9
1	A	471	ASP	3.9
2	B	70	LYS	3.9
1	A	135	ILE	3.8
2	B	431	LYS	3.7
2	B	94	ILE	3.6
1	A	69	THR	3.3
1	A	448	ARG	3.3

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Mol	Chain	Res	Type	RSRZ
1	A	447	ASN	3.3
1	A	468	THR	3.2
2	B	355	ALA	3.2
2	B	214	LEU	3.1
1	A	53	GLU	3.1
1	A	89	GLU	3.1
1	A	473	THR	3.0
2	B	426	TRP	3.0
2	B	95	PRO	3.0
1	A	138	GLU	2.9
2	B	195	ILE	2.9
1	A	358	ARG	2.8
2	B	5	ILE	2.8
2	B	7	THR	2.8
1	A	514	GLU	2.8
2	B	184	MET	2.8
1	A	272	PRO	2.7
2	B	68	SER	2.7
1	A	356	ARG	2.7
2	B	376	THR	2.7
1	A	54	ASN	2.6
1	A	451	LYS	2.6
2	B	216	THR	2.6
1	A	63	ILE	2.6
1	A	472	THR	2.6
1	A	2	ILE	2.4
1	A	112	GLY	2.4
2	B	200	THR	2.3
2	B	410	TRP	2.3
2	B	87	PHE	2.3
1	A	221	HIS	2.3
2	B	334	GLN	2.2
2	B	159	ILE	2.1
1	A	458	VAL	2.1
2	B	202	ILE	2.0

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

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
1	CSD	A	280	8/9	0.93	0.13	-	46,53,80,91	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(\AA^2)	Q<0.9
3	PO4	A	1301	5/5	0.88	0.27	3.74	133,136,139,146	0
3	PO4	A	1300	5/5	0.79	0.28	2.16	148,151,153,157	0
4	GWI	A	999	38/38	0.97	0.14	-0.27	28,50,91,104	0

6.5 Other polymers [i](#)

There are no such residues in this entry.