



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 09:09 AM UTC

PDB ID : 3IRX / pdb_00003irx
Title : Crystal Structure of HIV-1 reverse transcriptase (RT) in complex with the Non-nucleoside RT Inhibitor (E)-S-Methyl 5-(1-(3,7-Dimethyl-2-oxo-2,3-dihydrobenzo[d]oxazol-5-yl)-5-(5-methyl-1,3,4-oxadiazol-2-yl)pent-1-enyl)-2-methoxy-3-methylbenzothioate.
Authors : Ho, W.C.; Arnold, E.
Deposited on : 2009-08-24
Resolution : 2.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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](#) ①) 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
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

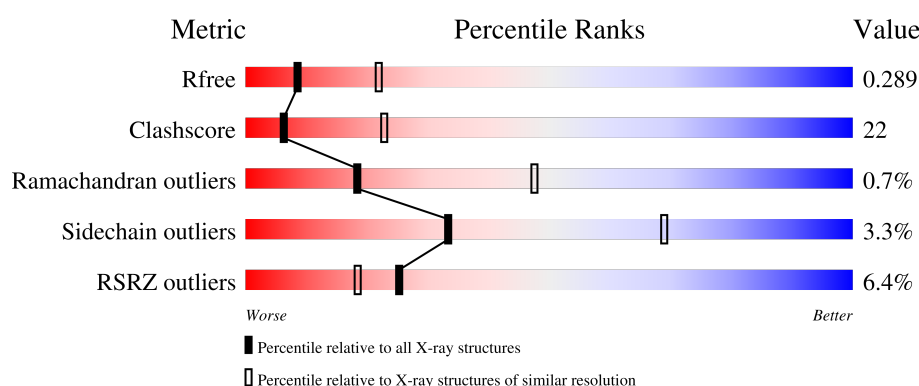
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	3866 (2.80-2.80)
Clashscore	190562	4276 (2.80-2.80)
Ramachandran outliers	187476	4196 (2.80-2.80)
Sidechain outliers	187428	4198 (2.80-2.80)
RSRZ outliers	180081	3869 (2.80-2.80)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments 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	558	7% 58% 38% ..
2	B	428	6% 61% 33% ..

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7985 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	555	Total	C	N	O	S	0	0	0
			4508	2918	749	834	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	GLY	-	expression tag	UNP P03366
A	-1	MET	-	expression tag	UNP P03366
A	0	VAL	-	expression tag	UNP P03366
A	172	ALA	LYS	engineered mutation	UNP P03366
A	173	ALA	LYS	engineered mutation	UNP P03366
A	280	SER	CYS	engineered mutation	UNP P03366

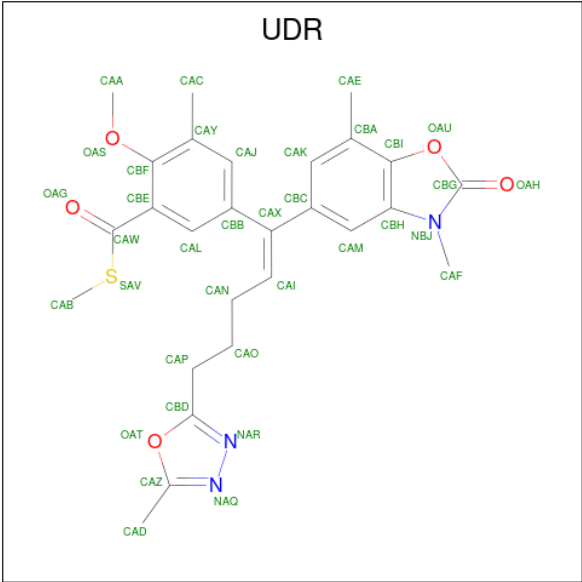
- Molecule 2 is a protein called Reverse transcriptase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	416	Total	C	N	O	S	0	0	0
			3441	2244	568	622	7			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	280	SER	CYS	engineered mutation	UNP P03366

- Molecule 3 is (E)-S-Methyl 5-(1-(3,7-Dimethyl-2-oxo-2,3-dihydrobenzo[d]oxazol-5-yl)-5-(5-methyl-1,3,4-oxadiazol-2-yl)pent-1-enyl)-2-methoxy-3-methyl benzothioate (CCD ID: UDR) (formula: C₂₇H₂₉N₃O₅S).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
3	A	1	36	27	3	5	1	0	0

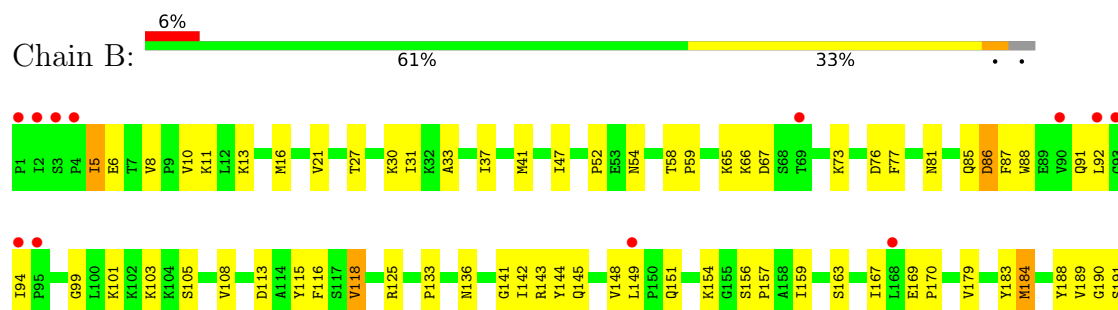
3 Residue-property plots [i](#)

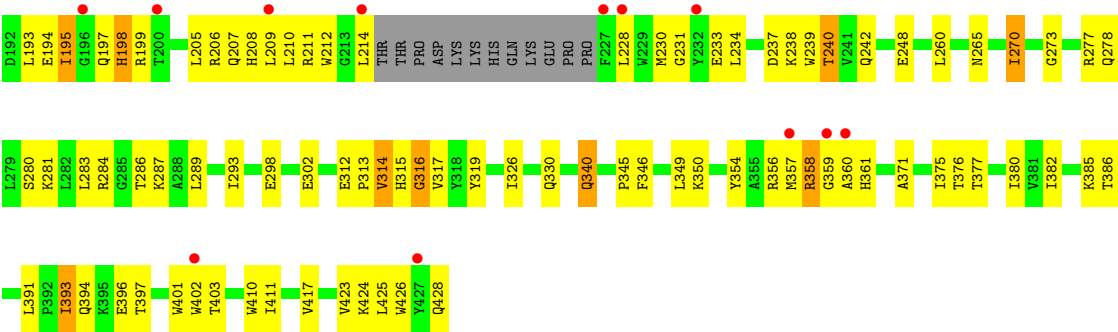
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: Reverse transcriptase/ribonuclease H



• Molecule 2: Reverse transcriptase





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	164.64Å 75.39Å 110.66Å 90.00° 100.08° 90.00°	Depositor
Resolution (Å)	50.00 – 2.80 50.00 – 2.80	Depositor EDS
% Data completeness (in resolution range)	87.1 (50.00-2.80) 87.1 (50.00-2.80)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.82 (at 2.81Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.236 , 0.289 0.237 , 0.289	Depositor DCC
R_{free} test set	1414 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å ²)	72.7	Xtriage
Anisotropy	0.096	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.28 , 52.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	7985	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.39% 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: UDR

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	0.57	1/4626 (0.0%)	0.96	17/6289 (0.3%)
2	B	0.55	0/3541	0.96	13/4810 (0.3%)
All	All	0.56	1/8167 (0.0%)	0.96	30/11099 (0.3%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	548	VAL	CA-CB	5.93	1.62	1.54

All (30) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	420	PRO	N-CA-C	8.90	121.56	110.70
1	A	545	ASN	N-CA-C	-8.82	101.34	112.90
2	B	240	THR	N-CA-C	-7.33	103.29	111.28
1	A	308	GLU	N-CA-C	-6.73	105.60	113.88
1	A	349	LEU	N-CA-C	-6.36	103.87	111.69
1	A	389	PHE	N-CA-C	6.24	120.22	110.17
2	B	349	LEU	N-CA-C	-6.19	105.67	113.72
2	B	237	ASP	N-CA-C	6.08	118.69	109.63
2	B	391	LEU	N-CA-C	5.78	117.17	109.83
1	A	442	VAL	N-CA-C	5.75	118.02	109.17
1	A	493	VAL	N-CA-C	5.67	116.52	108.42
2	B	289	LEU	N-CA-C	5.64	118.37	111.82
1	A	223	LYS	N-CA-C	5.58	118.72	109.46
2	B	273	GLY	N-CA-C	5.52	120.74	114.67
1	A	342	TYR	N-CA-C	5.49	115.78	108.38
2	B	184	MET	CB-CA-C	-5.48	109.78	117.23
1	A	352	GLY	N-CA-C	5.36	118.36	110.80
2	B	198	HIS	N-CA-C	-5.36	105.55	112.68

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	180	ILE	CB-CA-C	-5.33	104.38	110.73
1	A	508	ALA	N-CA-C	-5.28	106.68	113.23
2	B	118	VAL	CA-C-N	5.25	125.15	119.85
2	B	118	VAL	C-N-CA	5.25	125.15	119.85
2	B	189	VAL	N-CA-C	5.24	115.13	107.37
1	A	149	LEU	CA-C-N	5.17	125.09	119.76
1	A	149	LEU	C-N-CA	5.17	125.09	119.76
2	B	81	ASN	N-CA-C	5.12	116.55	111.07
1	A	322	SER	N-CA-C	5.09	116.83	111.28
1	A	391	LEU	N-CA-C	5.09	118.37	109.82
1	A	180	ILE	N-CA-C	5.09	115.78	107.24
2	B	141	GLY	N-CA-C	5.01	118.05	111.03

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	4508	0	4563	227	0
2	B	3441	0	3476	143	0
3	A	36	0	29	8	0
All	All	7985	0	8068	357	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 22.

All (357) 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:224:GLU:C	1:A:226:PRO:HD2	1.66	1.20
1:A:420:PRO:HB2	1:A:421:PRO:HD3	1.18	1.17
1:A:218:ASP:O	1:A:219:LYS:HD2	1.58	1.03
1:A:106:VAL:HG21	3:A:556:UDR:HAAB	1.42	0.99
2:B:315:HIS:O	2:B:317:VAL:HG23	1.65	0.96

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:420:PRO:HB2	1:A:421:PRO:CD	1.95	0.94
1:A:223:LYS:O	1:A:224:GLU:CD	2.10	0.94
1:A:421:PRO:O	1:A:422:LEU:HD23	1.68	0.94
1:A:106:VAL:HG21	3:A:556:UDR:CAA	1.98	0.93
1:A:34:LEU:HD21	1:A:62:ALA:HB2	1.51	0.92
1:A:399:GLU:HA	1:A:402:TRP:CD1	2.04	0.92
1:A:222:GLN:O	1:A:222:GLN:HG3	1.67	0.92
1:A:408:ALA:O	2:B:393:ILE:HD12	1.72	0.89
1:A:419:THR:HB	1:A:420:PRO:CD	2.03	0.88
1:A:103:LYS:HA	1:A:192:ASP:OD1	1.74	0.86
2:B:358:ARG:CZ	2:B:358:ARG:H	1.88	0.86
1:A:543:GLY:HA2	1:A:546:GLU:HB2	1.56	0.86
1:A:26:LEU:HB2	1:A:31:ILE:HD11	1.58	0.85
1:A:317:VAL:HG23	1:A:349:LEU:HD23	1.59	0.84
1:A:219:LYS:HE3	1:A:219:LYS:HA	1.60	0.83
1:A:223:LYS:O	1:A:224:GLU:OE1	1.95	0.82
2:B:37:ILE:HG22	2:B:41:MET:HE2	1.59	0.82
1:A:225:PRO:N	1:A:226:PRO:HD2	1.94	0.82
1:A:543:GLY:H	1:A:546:GLU:CG	1.93	0.82
1:A:224:GLU:C	1:A:226:PRO:CD	2.54	0.78
1:A:254:VAL:HB	1:A:290:THR:HA	1.64	0.78
1:A:543:GLY:H	1:A:546:GLU:HG2	1.49	0.78
2:B:66:LYS:HG3	2:B:67:ASP:H	1.49	0.78
2:B:170:PRO:HB2	2:B:208:HIS:CE1	2.19	0.77
2:B:33:ALA:O	2:B:37:ILE:HG12	1.85	0.76
1:A:225:PRO:N	1:A:226:PRO:CD	2.48	0.76
1:A:169:GLU:HB3	1:A:170:PRO:HD3	1.68	0.75
2:B:5:ILE:HD13	2:B:5:ILE:O	1.86	0.75
1:A:419:THR:HB	1:A:420:PRO:HD3	1.69	0.75
1:A:399:GLU:HG3	1:A:402:TRP:HE1	1.53	0.74
1:A:439:THR:H	1:A:460:ASN:ND2	1.85	0.74
1:A:419:THR:CB	1:A:420:PRO:CD	2.65	0.74
1:A:219:LYS:HA	1:A:219:LYS:CE	2.16	0.73
2:B:94:ILE:O	2:B:94:ILE:HD12	1.88	0.73
1:A:142:ILE:HD12	1:A:142:ILE:H	1.53	0.73
2:B:402:TRP:NE1	2:B:403:THR:HG23	2.03	0.73
2:B:31:ILE:HD11	2:B:133:PRO:HG2	1.71	0.72
1:A:516:GLU:O	1:A:520:GLN:HG3	1.90	0.71
1:A:539:HIS:O	1:A:540:LYS:HG2	1.91	0.71
1:A:541:GLY:HA2	1:A:546:GLU:HG3	1.72	0.71
1:A:114:ALA:HB1	1:A:214:LEU:HD22	1.72	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:458:VAL:HG23	2:B:286:THR:HG21	1.74	0.70
1:A:420:PRO:CB	1:A:421:PRO:HD3	2.09	0.70
1:A:142:ILE:HD12	1:A:142:ILE:N	2.06	0.70
2:B:402:TRP:CE2	2:B:403:THR:CG2	2.74	0.69
1:A:539:HIS:C	1:A:540:LYS:HG2	2.17	0.69
2:B:209:LEU:O	2:B:214:LEU:HB2	1.92	0.69
2:B:402:TRP:CZ2	2:B:403:THR:HG22	2.28	0.69
2:B:240:THR:O	2:B:350:LYS:HD2	1.93	0.68
1:A:399:GLU:HA	1:A:402:TRP:HD1	1.59	0.68
2:B:183:TYR:CD2	2:B:380:ILE:HD13	2.29	0.68
2:B:270:ILE:HD12	2:B:270:ILE:N	2.09	0.67
2:B:402:TRP:CE2	2:B:403:THR:HG23	2.29	0.67
1:A:419:THR:HB	1:A:420:PRO:HD2	1.77	0.66
1:A:540:LYS:O	1:A:546:GLU:CD	2.39	0.66
2:B:170:PRO:HB2	2:B:208:HIS:HE1	1.61	0.66
1:A:457:TYR:CD1	1:A:457:TYR:C	2.74	0.65
2:B:41:MET:HE1	2:B:73:LYS:CE	2.26	0.65
2:B:210:LEU:O	2:B:210:LEU:HD23	1.96	0.65
1:A:543:GLY:HA3	2:B:284:ARG:HA	1.79	0.65
1:A:458:VAL:CG2	2:B:286:THR:HG21	2.26	0.65
2:B:5:ILE:HG23	2:B:6:GLU:HG3	1.79	0.64
1:A:94:ILE:HD13	1:A:94:ILE:H	1.62	0.64
1:A:541:GLY:O	2:B:280:SER:HB3	1.99	0.63
2:B:423:VAL:HG12	2:B:426:TRP:CE3	2.33	0.63
1:A:164:MET:HG2	1:A:182:GLN:NE2	2.14	0.63
1:A:62:ALA:C	1:A:63:ILE:HD12	2.23	0.63
1:A:317:VAL:CG2	1:A:349:LEU:HD23	2.27	0.62
2:B:85:GLN:O	2:B:85:GLN:HG3	1.99	0.62
1:A:544:GLY:HA2	2:B:286:THR:CG2	2.30	0.62
2:B:10:VAL:HG13	2:B:85:GLN:HE21	1.64	0.62
1:A:16:MET:HE2	1:A:83:ARG:HG2	1.82	0.62
2:B:163:SER:O	2:B:167:ILE:HG13	2.00	0.62
1:A:233:GLU:OE1	1:A:243:PRO:HD3	1.99	0.61
1:A:402:TRP:CE2	1:A:403:THR:HG23	2.35	0.61
2:B:393:ILE:HG13	2:B:394:GLN:N	2.14	0.61
1:A:354:TYR:OH	1:A:370:GLU:HB3	2.00	0.61
1:A:202:ILE:O	1:A:206:ARG:HG3	2.00	0.60
2:B:108:VAL:O	2:B:231:GLY:HA2	2.01	0.60
1:A:181:TYR:CD2	3:A:556:UDR:HAPA	2.36	0.60
2:B:238:LYS:HG2	2:B:239:TRP:N	2.16	0.60
2:B:277:ARG:NH1	2:B:281:LYS:HZ2	2.00	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:399:GLU:HA	1:A:402:TRP:NE1	2.17	0.59
1:A:444:GLY:HA2	1:A:552:VAL:HG11	1.84	0.59
1:A:103:LYS:HD3	1:A:192:ASP:OD1	2.03	0.59
2:B:108:VAL:HG22	2:B:188:TYR:CD2	2.37	0.59
1:A:102:LYS:N	1:A:102:LYS:HE3	2.16	0.59
1:A:203:GLU:OE1	1:A:206:ARG:HD3	2.00	0.59
2:B:319:TYR:OH	2:B:385:LYS:HE2	2.02	0.59
1:A:220:LYS:O	1:A:221:HIS:ND1	2.35	0.59
2:B:402:TRP:CE2	2:B:403:THR:HG22	2.37	0.59
2:B:206:ARG:HB2	2:B:206:ARG:NH1	2.18	0.59
2:B:156:SER:N	2:B:157:PRO:HD2	2.18	0.59
1:A:257:ILE:O	1:A:261:VAL:HG23	2.02	0.59
1:A:361:HIS:NE2	1:A:505:ILE:HD12	2.17	0.59
1:A:305:GLU:O	1:A:309:ILE:HD13	2.02	0.59
1:A:65:LYS:HE3	1:A:68:SER:OG	2.02	0.58
1:A:419:THR:CB	1:A:420:PRO:HD2	2.32	0.58
1:A:323:LYS:HA	1:A:323:LYS:HE2	1.84	0.58
1:A:424:LYS:HE2	1:A:426:TRP:CZ3	2.39	0.58
2:B:113:ASP:HB2	2:B:228:LEU:HD21	1.84	0.58
2:B:423:VAL:HG12	2:B:426:TRP:CZ3	2.39	0.58
1:A:88:TRP:CD1	2:B:143:ARG:HD2	2.39	0.58
1:A:64:LYS:HB2	1:A:71:TRP:CD2	2.39	0.58
2:B:21:VAL:HB	2:B:59:PRO:HD3	1.85	0.58
2:B:358:ARG:CZ	2:B:358:ARG:N	2.64	0.58
1:A:417:VAL:O	1:A:417:VAL:HG13	2.04	0.57
2:B:76:ASP:OD1	2:B:76:ASP:C	2.47	0.57
1:A:28:GLU:HG3	1:A:135:ILE:HD12	1.86	0.57
1:A:544:GLY:HA2	2:B:286:THR:HG22	1.86	0.57
2:B:31:ILE:CD1	2:B:133:PRO:HG2	2.34	0.57
2:B:357:MET:O	2:B:357:MET:HG3	2.04	0.56
1:A:543:GLY:C	1:A:545:ASN:H	2.13	0.56
1:A:498:ASP:HB3	1:A:545:ASN:HD21	1.69	0.56
2:B:195:ILE:HD11	2:B:199:ARG:NH2	2.21	0.56
1:A:107:THR:HG22	1:A:108:VAL:N	2.20	0.56
1:A:266:TRP:O	1:A:269:GLN:HG2	2.06	0.56
1:A:278:GLN:O	1:A:282:LEU:HG	2.06	0.55
1:A:246:LEU:HD21	1:A:310:LEU:HD12	1.87	0.55
1:A:366:LYS:HE2	1:A:405:TYR:OH	2.06	0.55
1:A:543:GLY:CA	1:A:546:GLU:HB2	2.33	0.55
1:A:542:ILE:O	2:B:283:LEU:HB3	2.08	0.55
2:B:402:TRP:CG	2:B:403:THR:N	2.75	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:246:LEU:HD21	1:A:310:LEU:CD1	2.37	0.54
1:A:451:LYS:HB3	1:A:471:ASN:HA	1.89	0.54
2:B:376:THR:CG2	2:B:386:THR:HG22	2.37	0.54
1:A:17:ASP:O	1:A:83:ARG:HD3	2.07	0.54
1:A:10:VAL:HG12	1:A:11:LYS:N	2.21	0.54
1:A:222:GLN:O	1:A:222:GLN:CG	2.48	0.54
1:A:363:ASN:HD21	1:A:366:LYS:HG3	1.73	0.54
1:A:219:LYS:O	1:A:220:LYS:HB2	2.07	0.54
2:B:356:ARG:O	2:B:356:ARG:HG2	2.08	0.54
1:A:39:THR:O	1:A:43:LYS:HG2	2.08	0.54
2:B:233:GLU:CD	2:B:233:GLU:H	2.16	0.53
1:A:498:ASP:HB3	1:A:545:ASN:ND2	2.23	0.53
2:B:315:HIS:O	2:B:317:VAL:CG2	2.50	0.53
2:B:118:VAL:HB	2:B:149:LEU:HD12	1.89	0.53
2:B:315:HIS:O	2:B:316:GLY:C	2.52	0.53
1:A:345:PRO:O	1:A:346:PHE:HB2	2.09	0.53
1:A:31:ILE:O	1:A:35:VAL:HG23	2.09	0.52
2:B:91:GLN:O	2:B:92:LEU:HB2	2.09	0.52
1:A:29:GLU:HG3	1:A:30:LYS:N	2.24	0.52
1:A:376:THR:O	1:A:380:ILE:HG12	2.10	0.52
1:A:420:PRO:CB	1:A:421:PRO:CD	2.71	0.52
2:B:193:LEU:HB3	2:B:197:GLN:HG3	1.92	0.52
1:A:398:TRP:CZ2	1:A:411:ILE:HG12	2.44	0.52
1:A:183:TYR:O	1:A:184:MET:HB2	2.10	0.52
1:A:258:GLN:HB2	1:A:283:LEU:HD22	1.92	0.52
1:A:284:ARG:H	1:A:284:ARG:HD3	1.75	0.52
2:B:115:TYR:OH	2:B:157:PRO:HA	2.10	0.52
2:B:183:TYR:CE2	2:B:184:MET:HG2	2.45	0.52
2:B:393:ILE:HD11	2:B:397:THR:CG2	2.39	0.52
2:B:85:GLN:O	2:B:87:PHE:N	2.43	0.52
2:B:376:THR:HG23	2:B:386:THR:HG22	1.90	0.52
1:A:175:ASN:N	1:A:176:PRO:HD3	2.25	0.51
1:A:536:VAL:HG21	1:A:542:ILE:HG21	1.92	0.51
1:A:65:LYS:HE3	1:A:68:SER:CB	2.39	0.51
1:A:363:ASN:HA	1:A:511:ASP:OD1	2.10	0.51
2:B:169:GLU:HB2	2:B:170:PRO:HD3	1.91	0.51
2:B:359:GLY:C	2:B:361:HIS:H	2.18	0.51
1:A:53:GLU:O	1:A:55:PRO:HD3	2.11	0.51
1:A:277:ARG:NH1	1:A:334:GLN:O	2.44	0.51
1:A:449:GLU:N	1:A:449:GLU:CD	2.69	0.51
1:A:540:LYS:HE3	2:B:265:ASN:OD1	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:183:TYR:CD2	2:B:184:MET:HG2	2.45	0.51
1:A:470:THR:O	1:A:471:ASN:HB3	2.11	0.50
1:A:483:TYR:CE2	1:A:487:GLN:NE2	2.80	0.50
1:A:311:LYS:O	1:A:313:PRO:HD3	2.11	0.50
1:A:457:TYR:C	1:A:457:TYR:HD1	2.19	0.50
2:B:270:ILE:N	2:B:270:ILE:CD1	2.75	0.50
2:B:345:PRO:O	2:B:346:PHE:HB2	2.11	0.50
1:A:317:VAL:HG23	1:A:349:LEU:CD2	2.38	0.50
1:A:155:GLY:O	1:A:159:ILE:HG12	2.10	0.50
2:B:41:MET:HE1	2:B:73:LYS:HE3	1.93	0.50
1:A:544:GLY:O	1:A:548:VAL:HG12	2.10	0.50
2:B:125:ARG:O	2:B:145:GLN:HG3	2.12	0.50
1:A:183:TYR:CE2	1:A:230:MET:HE3	2.47	0.50
2:B:423:VAL:C	2:B:425:LEU:H	2.19	0.50
1:A:106:VAL:HG21	3:A:556:UDR:HAAA	1.90	0.50
2:B:410:TRP:O	2:B:411:ILE:HD13	2.12	0.49
1:A:207:GLN:O	1:A:210:LEU:HB3	2.11	0.49
1:A:543:GLY:C	1:A:545:ASN:N	2.70	0.49
1:A:106:VAL:CG2	3:A:556:UDR:HAAB	2.27	0.49
2:B:425:LEU:HD12	2:B:428:GLN:HB2	1.94	0.49
1:A:245:VAL:O	1:A:247:PRO:HD3	2.10	0.49
2:B:101:LYS:HD3	2:B:382:ILE:HG23	1.94	0.49
2:B:142:ILE:CG2	2:B:144:TYR:CE1	2.96	0.49
1:A:65:LYS:O	1:A:68:SER:HB2	2.13	0.49
1:A:74:LEU:HD13	1:A:74:LEU:C	2.38	0.49
1:A:443:ASP:HB2	1:A:548:VAL:HG13	1.94	0.49
1:A:31:ILE:HD12	1:A:133:PRO:HG2	1.95	0.49
1:A:326:ILE:HG22	1:A:327:ALA:N	2.28	0.49
2:B:10:VAL:O	2:B:11:LYS:HD3	2.13	0.49
1:A:129:ALA:HA	1:A:144:TYR:O	2.13	0.49
1:A:224:GLU:HB3	1:A:226:PRO:HD2	1.94	0.49
1:A:224:GLU:O	1:A:226:PRO:HD2	2.07	0.49
2:B:278:GLN:HB2	2:B:302:GLU:OE1	2.13	0.49
1:A:223:LYS:C	1:A:224:GLU:CD	2.80	0.48
1:A:541:GLY:O	2:B:280:SER:O	2.29	0.48
2:B:281:LYS:HD2	2:B:284:ARG:NH1	2.28	0.48
1:A:5:ILE:O	1:A:6:GLU:C	2.56	0.48
2:B:206:ARG:HB2	2:B:206:ARG:HH11	1.77	0.48
1:A:78:ARG:HH11	1:A:78:ARG:HB2	1.78	0.48
1:A:181:TYR:CE2	1:A:183:TYR:HB2	2.48	0.48
2:B:312:GLU:HB3	2:B:313:PRO:HD2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:433:PRO:HA	1:A:532:TYR:CD2	2.49	0.48
2:B:287:LYS:HD3	2:B:293:ILE:HD11	1.96	0.48
1:A:63:ILE:HD12	1:A:63:ILE:N	2.29	0.47
1:A:229:TRP:CE2	3:A:556:UDR:HAO	2.50	0.47
1:A:224:GLU:O	1:A:227:PHE:CE2	2.67	0.47
2:B:376:THR:HG21	2:B:410:TRP:CZ3	2.49	0.47
1:A:426:TRP:HD1	1:A:511:ASP:HB2	1.80	0.47
1:A:47:ILE:HD12	1:A:144:TYR:CD2	2.50	0.47
1:A:142:ILE:H	1:A:142:ILE:CD1	2.23	0.47
2:B:167:ILE:HG12	2:B:212:TRP:CE3	2.49	0.47
1:A:244:ILE:HG13	1:A:267:ALA:HB2	1.97	0.47
2:B:195:ILE:HD11	2:B:199:ARG:HH21	1.80	0.47
1:A:284:ARG:HD3	1:A:284:ARG:N	2.30	0.47
1:A:356:ARG:HG2	1:A:357:MET:O	2.15	0.47
1:A:536:VAL:HB	1:A:542:ILE:HG13	1.96	0.47
1:A:295:LEU:HD12	1:A:295:LEU:N	2.30	0.46
1:A:439:THR:H	1:A:460:ASN:HD22	1.61	0.46
1:A:31:ILE:CD1	1:A:133:PRO:HG2	2.44	0.46
1:A:544:GLY:CA	2:B:286:THR:HG22	2.44	0.46
1:A:167:ILE:O	1:A:170:PRO:HD2	2.15	0.46
2:B:47:ILE:HD12	2:B:144:TYR:CD2	2.50	0.46
2:B:314:VAL:HG22	2:B:315:HIS:N	2.31	0.46
3:A:556:UDR:HANA	3:A:556:UDR:CAL	2.44	0.46
2:B:8:VAL:O	2:B:10:VAL:HG23	2.16	0.46
1:A:246:LEU:CD1	1:A:307:ARG:HG2	2.45	0.46
1:A:270:ILE:O	1:A:272:PRO:HD3	2.16	0.46
1:A:293:ILE:HG23	1:A:294:PRO:HD2	1.97	0.46
2:B:208:HIS:O	2:B:211:ARG:HG2	2.16	0.46
1:A:148:VAL:O	1:A:150:PRO:HD3	2.16	0.46
2:B:65:LYS:HB2	2:B:230:MET:HE1	1.97	0.46
1:A:102:LYS:C	1:A:102:LYS:HD2	2.41	0.45
1:A:424:LYS:HE2	1:A:426:TRP:CE3	2.51	0.45
2:B:116:PHE:CD2	2:B:148:VAL:HG21	2.51	0.45
1:A:426:TRP:CD1	1:A:511:ASP:HB2	2.51	0.45
2:B:5:ILE:HD13	2:B:5:ILE:C	2.41	0.45
1:A:46:LYS:N	1:A:46:LYS:HE3	2.32	0.45
1:A:363:ASN:ND2	1:A:366:LYS:HG3	2.31	0.45
2:B:105:SER:O	2:B:190:GLY:HA2	2.16	0.45
2:B:115:TYR:HE2	2:B:156:SER:O	1.99	0.45
2:B:207:GLN:OE1	2:B:210:LEU:HD13	2.17	0.45
1:A:101:LYS:HG3	1:A:103:LYS:NZ	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:250:ASP:OD2	1:A:251:SER:N	2.50	0.45
2:B:8:VAL:HB	2:B:159:ILE:HD12	1.98	0.45
2:B:326:ILE:N	2:B:326:ILE:HD12	2.31	0.45
1:A:183:TYR:HE2	1:A:230:MET:HE3	1.80	0.45
1:A:287:LYS:O	1:A:288:ALA:HB3	2.15	0.45
2:B:340:GLN:HE21	2:B:340:GLN:HB2	1.60	0.45
1:A:224:GLU:CB	1:A:226:PRO:HD2	2.47	0.45
1:A:356:ARG:HD3	1:A:367:GLN:NE2	2.32	0.45
2:B:371:ALA:O	2:B:375:ILE:HG13	2.17	0.45
1:A:19:PRO:O	1:A:20:LYS:HD3	2.17	0.45
1:A:120:LEU:HD12	1:A:121:ASP:H	1.82	0.45
1:A:229:TRP:O	1:A:230:MET:C	2.60	0.45
1:A:474:ASN:O	1:A:478:GLU:HG3	2.16	0.45
2:B:41:MET:HE1	2:B:73:LYS:HE2	1.96	0.45
2:B:169:GLU:OE2	2:B:169:GLU:HA	2.16	0.45
1:A:417:VAL:O	1:A:417:VAL:CG1	2.65	0.44
1:A:199:ARG:HB2	1:A:199:ARG:CZ	2.47	0.44
2:B:11:LYS:H	2:B:85:GLN:NE2	2.15	0.44
2:B:85:GLN:O	2:B:86:ASP:C	2.61	0.44
1:A:118:VAL:HG11	1:A:149:LEU:HD11	1.98	0.44
1:A:542:ILE:O	1:A:542:ILE:HG22	2.16	0.44
2:B:169:GLU:N	2:B:170:PRO:CD	2.81	0.44
2:B:359:GLY:C	2:B:361:HIS:N	2.76	0.44
1:A:70:LYS:HG2	1:A:71:TRP:H	1.82	0.44
1:A:553:SER:O	1:A:554:ALA:C	2.60	0.44
1:A:42:GLU:OE1	1:A:49:LYS:HG3	2.18	0.44
1:A:64:LYS:HB2	1:A:71:TRP:CE2	2.53	0.44
1:A:281:LYS:O	1:A:281:LYS:HD3	2.17	0.44
2:B:66:LYS:HG3	2:B:67:ASP:N	2.24	0.44
2:B:234:LEU:HD12	2:B:234:LEU:O	2.18	0.44
1:A:398:TRP:CH2	1:A:411:ILE:HG12	2.53	0.44
1:A:194:GLU:HG3	1:A:197:GLN:NE2	2.33	0.44
2:B:108:VAL:HG22	2:B:188:TYR:CE2	2.53	0.44
2:B:358:ARG:N	2:B:358:ARG:NE	2.66	0.44
2:B:402:TRP:CZ2	2:B:403:THR:CG2	2.98	0.43
1:A:193:LEU:O	1:A:194:GLU:C	2.60	0.43
1:A:194:GLU:OE1	1:A:194:GLU:HA	2.18	0.43
2:B:191:SER:OG	2:B:198:HIS:ND1	2.44	0.43
1:A:0:VAL:N	1:A:1:PRO:CD	2.82	0.43
1:A:102:LYS:HD2	1:A:102:LYS:O	2.18	0.43
1:A:156:SER:N	1:A:157:PRO:CD	2.82	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:314:VAL:CG2	2:B:315:HIS:N	2.81	0.43
1:A:254:VAL:HG23	1:A:291:GLU:N	2.33	0.43
2:B:85:GLN:O	2:B:85:GLN:CG	2.65	0.43
2:B:151:GLN:HE21	2:B:151:GLN:HB3	1.61	0.43
1:A:99:GLY:HA3	2:B:136:ASN:ND2	2.32	0.43
2:B:27:THR:OG1	2:B:30:LYS:HG3	2.18	0.43
2:B:154:LYS:HG2	2:B:154:LYS:O	2.18	0.43
2:B:354:TYR:OH	2:B:357:MET:HG2	2.19	0.43
1:A:449:GLU:N	1:A:449:GLU:OE2	2.51	0.43
1:A:498:ASP:C	1:A:498:ASP:OD1	2.62	0.43
2:B:116:PHE:HD2	2:B:148:VAL:HG21	1.83	0.42
2:B:396:GLU:CD	2:B:396:GLU:H	2.27	0.42
1:A:438:GLU:HG3	1:A:461:LYS:HD3	2.01	0.42
1:A:115:TYR:OH	1:A:156:SER:HB3	2.19	0.42
1:A:454:LYS:HB2	1:A:552:VAL:HG13	2.02	0.42
2:B:239:TRP:O	2:B:242:GLN:NE2	2.52	0.42
1:A:53:GLU:H	1:A:53:GLU:CD	2.26	0.42
1:A:190:GLY:O	1:A:191:SER:HB3	2.19	0.42
1:A:96:HIS:O	1:A:97:PRO:C	2.62	0.42
2:B:358:ARG:HA	2:B:358:ARG:NH1	2.34	0.42
1:A:543:GLY:N	1:A:546:GLU:HG2	2.27	0.42
2:B:52:PRO:C	2:B:54:ASN:N	2.76	0.42
1:A:194:GLU:HG3	1:A:197:GLN:HE21	1.85	0.42
1:A:284:ARG:H	1:A:284:ARG:CD	2.33	0.42
1:A:360:ALA:O	1:A:514:GLU:HG2	2.20	0.42
2:B:194:GLU:HG2	2:B:197:GLN:HG2	2.02	0.42
1:A:107:THR:HG22	1:A:108:VAL:H	1.85	0.41
1:A:246:LEU:HD11	1:A:310:LEU:HD12	2.02	0.41
1:A:354:TYR:CZ	1:A:370:GLU:HB3	2.54	0.41
2:B:401:TRP:O	2:B:402:TRP:C	2.63	0.41
1:A:37:ILE:O	1:A:40:GLU:HG2	2.21	0.41
1:A:70:LYS:HG2	1:A:71:TRP:N	2.35	0.41
2:B:13:LYS:HD2	2:B:16:MET:CE	2.50	0.41
1:A:218:ASP:O	1:A:219:LYS:CD	2.48	0.41
2:B:52:PRO:C	2:B:54:ASN:H	2.28	0.41
1:A:542:ILE:H	1:A:546:GLU:CG	2.34	0.41
2:B:330:GLN:HE21	2:B:330:GLN:HB2	1.70	0.41
1:A:206:ARG:CZ	1:A:218:ASP:CB	2.99	0.41
1:A:399:GLU:CA	1:A:402:TRP:CD1	2.92	0.41
1:A:29:GLU:CG	1:A:30:LYS:N	2.84	0.41
1:A:287:LYS:HB3	1:A:287:LYS:NZ	2.36	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:503:LEU:HA	1:A:503:LEU:HD23	1.64	0.41
1:A:541:GLY:C	2:B:280:SER:HB3	2.46	0.41
2:B:13:LYS:HD2	2:B:16:MET:HE1	2.02	0.41
2:B:99:GLY:O	2:B:103:LYS:HG2	2.21	0.41
2:B:205:LEU:O	2:B:209:LEU:HG	2.20	0.41
1:A:118:VAL:HA	1:A:119:PRO:HD2	1.95	0.41
2:B:156:SER:N	2:B:157:PRO:CD	2.84	0.41
1:A:283:LEU:HA	1:A:286:THR:HG23	2.03	0.40
2:B:260:LEU:HA	2:B:260:LEU:HD12	1.83	0.40
1:A:181:TYR:CD2	3:A:556:UDR:CAP	3.05	0.40
1:A:223:LYS:HE2	1:A:223:LYS:HB3	1.72	0.40
1:A:293:ILE:CG2	1:A:294:PRO:HD2	2.52	0.40
1:A:306:ASN:O	1:A:310:LEU:HB2	2.21	0.40
2:B:58:THR:HG21	2:B:77:PHE:CD2	2.57	0.40
2:B:103:LYS:HE2	2:B:179:VAL:HG23	2.04	0.40
2:B:360:ALA:O	2:B:361:HIS:CG	2.74	0.40
1:A:12:LEU:HD22	1:A:83:ARG:O	2.21	0.40
1:A:223:LYS:HB3	1:A:223:LYS:NZ	2.30	0.40
1:A:283:LEU:HD12	1:A:283:LEU:C	2.46	0.40
1:A:542:ILE:H	1:A:546:GLU:HG2	1.86	0.40
2:B:393:ILE:HD11	2:B:397:THR:HG21	2.03	0.40
1:A:175:ASN:HB3	1:A:178:ILE:HG12	2.03	0.40
1:A:391:LEU:O	1:A:417:VAL:HG12	2.22	0.40
2:B:298:GLU:CD	2:B:298:GLU:H	2.29	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	553/558 (99%)	482 (87%)	68 (12%)	3 (0%)	24 55

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	412/428 (96%)	362 (88%)	46 (11%)	4 (1%)	12	38
All	All	965/986 (98%)	844 (88%)	114 (12%)	7 (1%)	18	47

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	424	LYS
1	A	194	GLU
1	A	184	MET
2	B	86	ASP
2	B	316	GLY
1	A	317	VAL
2	B	195	ILE

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	494/495 (100%)	475 (96%)	19 (4%)	29	64
2	B	378/390 (97%)	368 (97%)	10 (3%)	40	75
All	All	872/885 (98%)	843 (97%)	29 (3%)	33	69

All (29) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	SER
1	A	46	LYS
1	A	78	ARG
1	A	91	GLN
1	A	92	LEU
1	A	94	ILE
1	A	102	LYS
1	A	142	ILE
1	A	192	ASP

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Mol	Chain	Res	Type
1	A	207	GLN
1	A	219	LYS
1	A	223	LYS
1	A	281	LYS
1	A	287	LYS
1	A	347	LYS
1	A	362	THR
1	A	459	THR
1	A	512	LYS
1	A	545	ASN
2	B	5	ILE
2	B	88	TRP
2	B	248	GLU
2	B	270	ILE
2	B	314	VAL
2	B	340	GLN
2	B	358	ARG
2	B	377	THR
2	B	393	ILE
2	B	417	VAL

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (28) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	23	GLN
1	A	85	GLN
1	A	151	GLN
1	A	197	GLN
1	A	222	GLN
1	A	242	GLN
1	A	255	ASN
1	A	306	ASN
1	A	315	HIS
1	A	336	GLN
1	A	394	GLN
1	A	407	GLN
1	A	460	ASN
1	A	475	GLN
1	A	487	GLN
1	A	507	GLN
1	A	524	GLN
1	A	545	ASN

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Mol	Chain	Res	Type
2	B	85	GLN
2	B	151	GLN
2	B	174	GLN
2	B	197	GLN
2	B	208	HIS
2	B	255	ASN
2	B	278	GLN
2	B	330	GLN
2	B	340	GLN
2	B	418	ASN

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 ⓘ

1 ligand 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	$\# Z > 2$
3	UDR	A	556	-	38,39,39	2.76	10 (26%)	50,56,56	3.20	16 (32%)

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	UDR	A	556	-	-	3/22/23/23	0/4/4/4

All (10) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	556	UDR	NAR-NAQ	-7.44	1.23	1.39
3	A	556	UDR	CBB-CAX	-7.39	1.38	1.49
3	A	556	UDR	CBC-CAX	-5.97	1.40	1.49
3	A	556	UDR	CAZ-NAQ	5.61	1.33	1.28
3	A	556	UDR	CAI-CAX	4.93	1.39	1.34
3	A	556	UDR	CBD-NAR	4.57	1.34	1.28
3	A	556	UDR	CAD-CAZ	3.73	1.55	1.48
3	A	556	UDR	CAN-CAI	-3.56	1.39	1.50
3	A	556	UDR	CBH-NBJ	-3.15	1.33	1.39
3	A	556	UDR	CBE-CAW	2.59	1.53	1.49

All (16) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	556	UDR	CAD-CAZ-NAQ	-11.15	123.67	129.21
3	A	556	UDR	CAB-SAV-CAW	9.31	108.45	99.75
3	A	556	UDR	OAU-CBI-CBA	7.85	132.33	124.59
3	A	556	UDR	CBD-NAR-NAQ	6.90	109.99	106.25
3	A	556	UDR	OAT-CAZ-NAQ	-5.69	110.21	112.34
3	A	556	UDR	CAZ-NAQ-NAR	5.56	108.66	106.42
3	A	556	UDR	OAT-CBD-NAR	-5.34	108.47	112.75
3	A	556	UDR	OAT-CAZ-CAD	4.30	126.11	117.49
3	A	556	UDR	OAU-CBG-OAH	3.31	125.50	122.43
3	A	556	UDR	CAP-CBD-NAR	-3.11	126.96	129.47
3	A	556	UDR	OAT-CBD-CAP	2.86	124.43	117.96
3	A	556	UDR	CBE-CAW-SAV	2.56	119.82	114.78
3	A	556	UDR	CBI-OAU-CBG	-2.23	105.96	107.18
3	A	556	UDR	CBC-CAX-CBB	2.19	119.65	116.92
3	A	556	UDR	CAA-OAS-CBF	-2.06	109.14	114.74
3	A	556	UDR	CBC-CAX-CAI	-2.00	119.23	121.76

There are no chirality outliers.

All (3) torsion outliers are listed below:

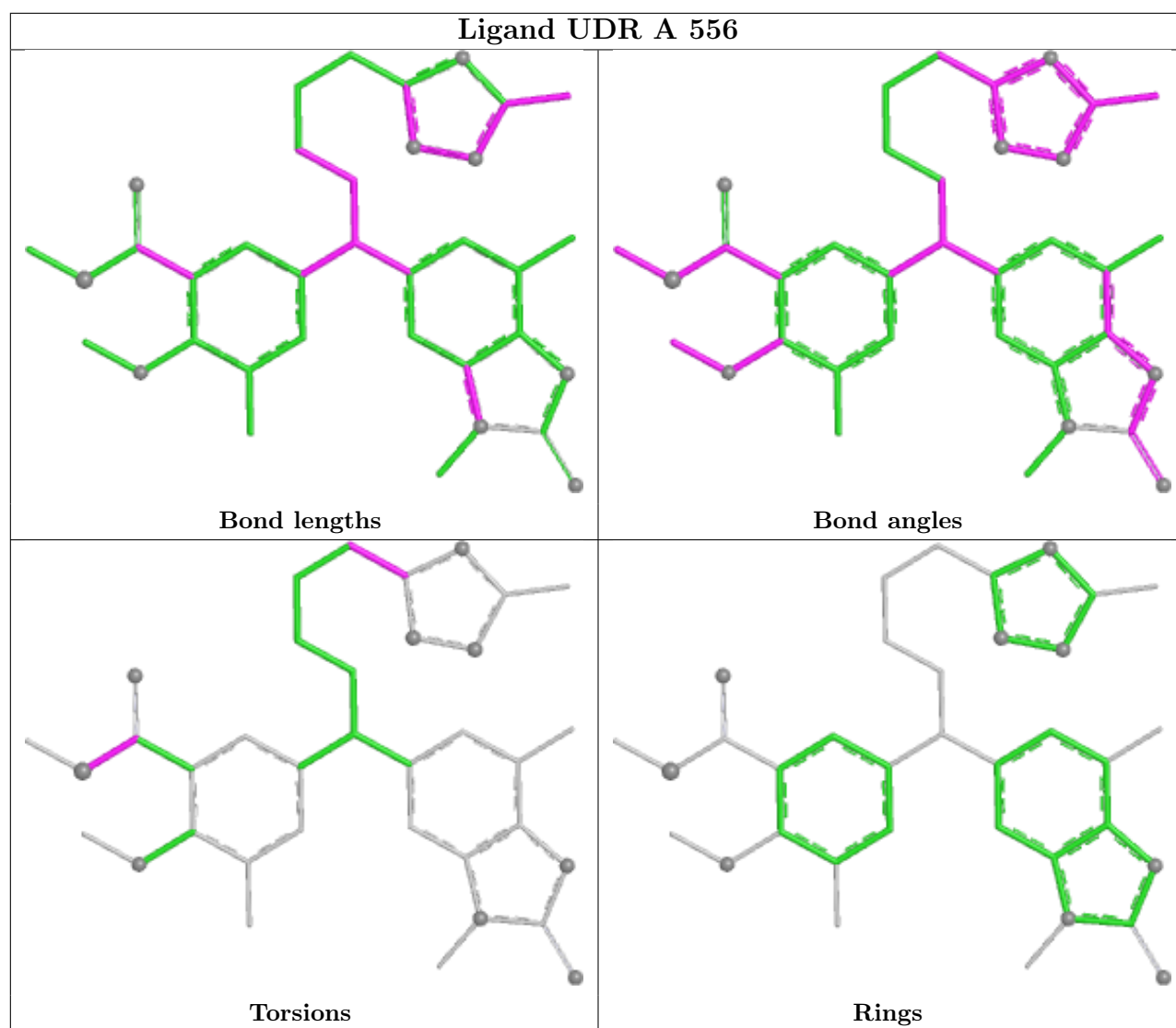
Mol	Chain	Res	Type	Atoms
3	A	556	UDR	CAO-CAP-CBD-NAR
3	A	556	UDR	OAG-CAW-SAV-CAB
3	A	556	UDR	CAO-CAP-CBD-OAT

There are no ring outliers.

1 monomer is involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	556	UDR	8	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	555/558 (99%)	0.48	38 (6%)	23 17	40, 90, 158, 175	0
2	B	416/428 (97%)	0.39	24 (5%)	29 22	49, 81, 149, 182	0
All	All	971/986 (98%)	0.44	62 (6%)	25 18	40, 86, 154, 182	0

All (62) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	545	ASN	6.7
1	A	548	VAL	6.1
1	A	547	GLN	5.9
1	A	549	ASP	5.7
1	A	541	GLY	5.7
1	A	551	LEU	5.2
1	A	552	VAL	5.1
1	A	553	SER	4.8
1	A	544	GLY	4.8
2	B	2	ILE	4.8
1	A	542	ILE	3.9
2	B	92	LEU	3.8
2	B	90	VAL	3.7
1	A	546	GLU	3.5
1	A	247	PRO	3.4
1	A	92	LEU	3.3
2	B	94	ILE	3.3
1	A	251	SER	3.2
2	B	228	LEU	3.1
1	A	550	LYS	3.1
2	B	93	GLY	3.0
2	B	402	TRP	2.9
2	B	214	LEU	2.9
2	B	95	PRO	2.9

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Mol	Chain	Res	Type	RSRZ
2	B	149	LEU	2.9
1	A	292	VAL	2.9
1	A	252	TRP	2.8
2	B	168	LEU	2.7
1	A	67	ASP	2.6
1	A	543	GLY	2.6
2	B	3	SER	2.5
2	B	4	PRO	2.5
1	A	257	ILE	2.5
1	A	24	TRP	2.5
1	A	283	LEU	2.5
1	A	317	VAL	2.5
1	A	303	LEU	2.4
1	A	195	ILE	2.4
2	B	232	TYR	2.4
2	B	200	THR	2.4
2	B	209	LEU	2.4
2	B	357	MET	2.4
2	B	359	GLY	2.3
1	A	22	LYS	2.3
2	B	360	ALA	2.3
1	A	244	ILE	2.3
1	A	299	ALA	2.3
1	A	554	ALA	2.3
1	A	132	ILE	2.3
1	A	90	VAL	2.2
2	B	227	PHE	2.2
2	B	69	THR	2.2
2	B	427	TYR	2.2
1	A	245	VAL	2.2
1	A	420	PRO	2.1
2	B	196	GLY	2.1
1	A	38	CYS	2.1
2	B	1	PRO	2.0
1	A	214	LEU	2.0
1	A	218	ASP	2.0
1	A	309	ILE	2.0
1	A	402	TRP	2.0

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

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates [i](#)

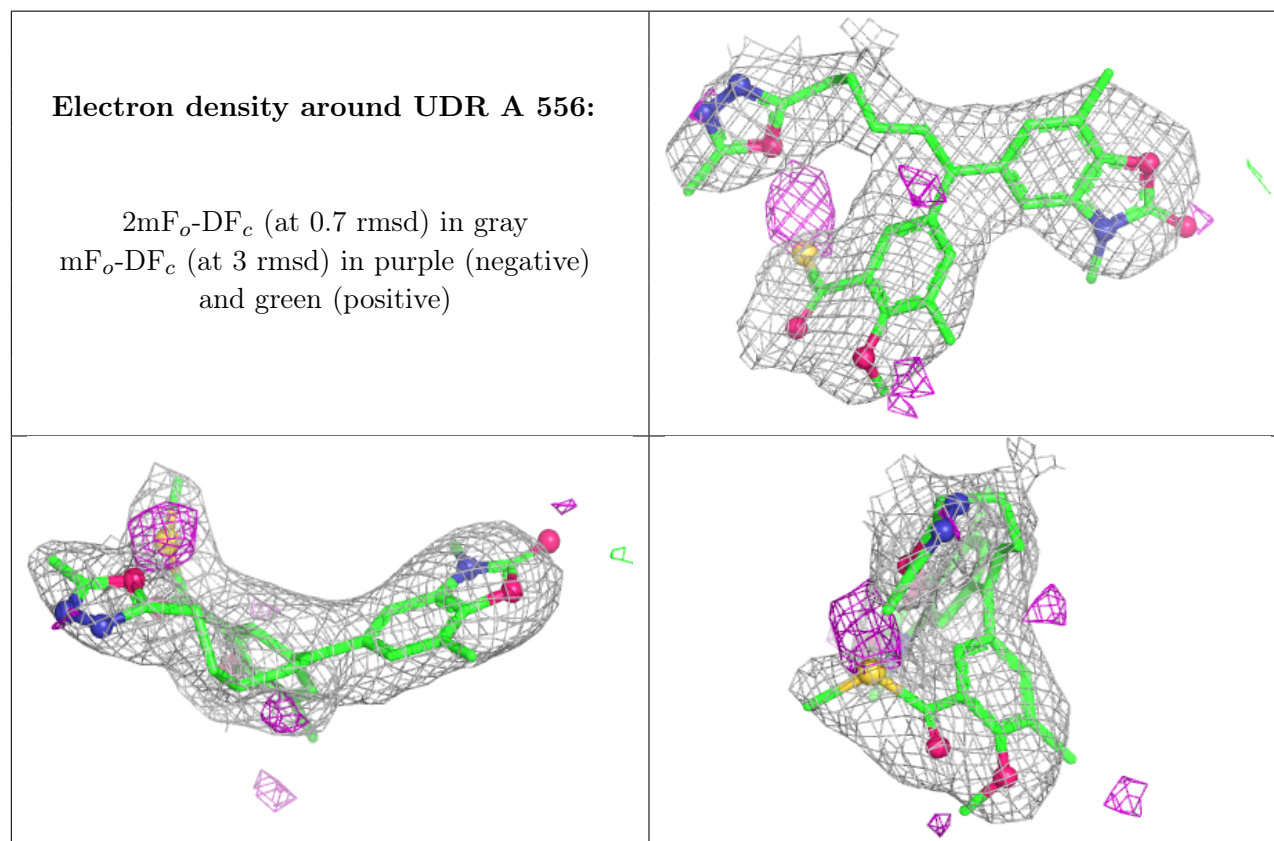
There are no oligosaccharides in this entry.

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	UDR	A	556	36/36	0.85	0.14	69,76,79,81	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.



6.5 Other polymers [i](#)

There are no such residues in this entry.