



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 07:39 PM UTC

PDB ID : 7S7V / pdb_00007s7v
Title : Crystal structure of iNicSnFR3a Fluorescent Nicotine Sensor
Authors : Fan, C.; Shivange, A.V.; Looger, L.L.; Lester, H.A.; Rees, D.C.
Deposited on : 2021-09-17
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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

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

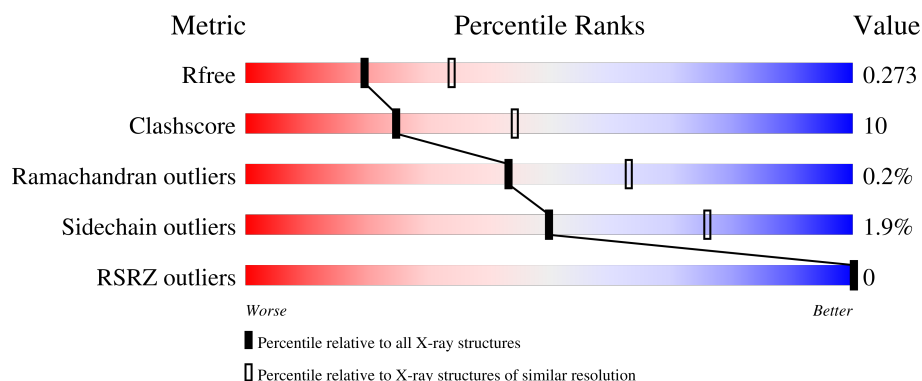
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

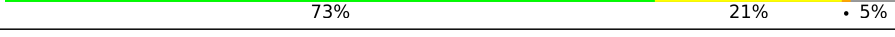
The reported resolution of this entry is 2.50 Å.

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}	180053	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (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 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	537	 72% 22% • 5%
1	B	537	 69% 24% • 5%
1	C	537	 73% 20% • 6%
1	D	537	 73% 21% • 5%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 16649 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 iNicSnFR 3.0 Fluorescent Nicotine Sensor.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	509	Total	C	N	O	S	0	0	0
			4040	2588	671	768	13			
1	B	509	Total	C	N	O	S	0	0	0
			4035	2584	670	768	13			
1	C	507	Total	C	N	O	S	0	0	0
			4024	2578	668	765	13			
1	D	509	Total	C	N	O	S	0	1	0
			4038	2586	670	769	13			

- Molecule 2 is water.

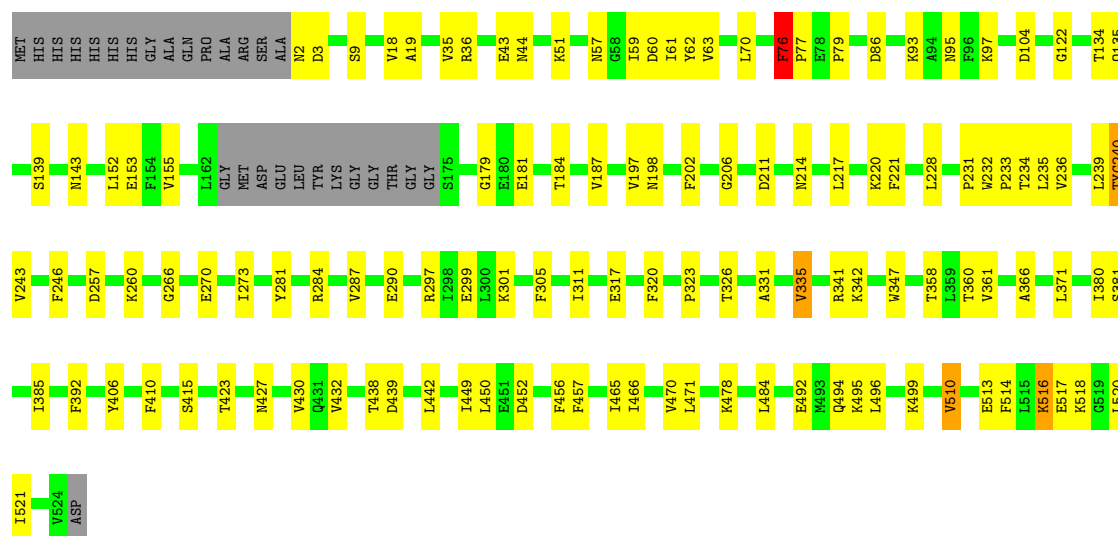
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	134	Total	O	0	0
			134	134		
2	B	112	Total	O	0	0
			112	112		
2	C	147	Total	O	0	0
			147	147		
2	D	119	Total	O	0	0
			119	119		

3 Residue-property plots

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

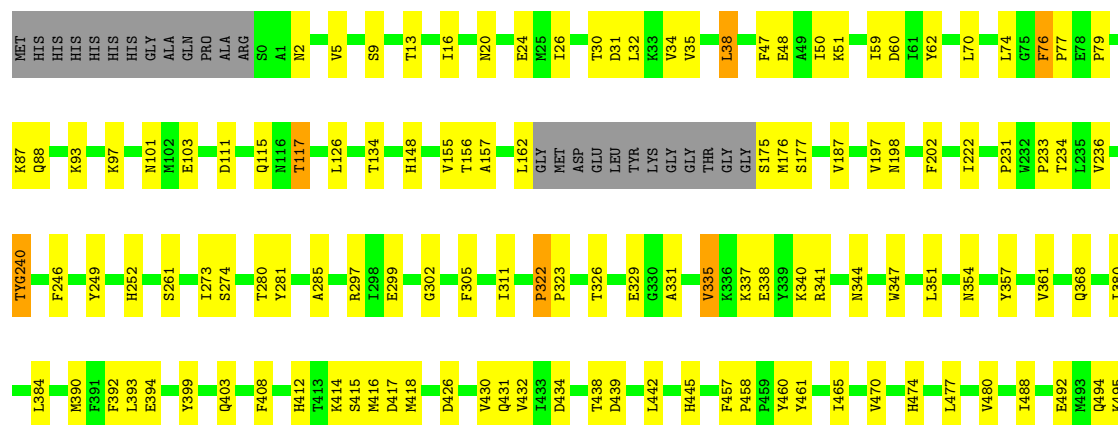
• Molecule 1: iNicSnFR 3.0 Fluorescent Nicotine Sensor

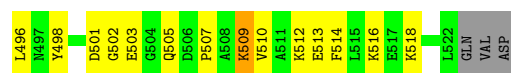
Chain A: 



• Molecule 1: iNicSnFR 3.0 Fluorescent Nicotine Sensor

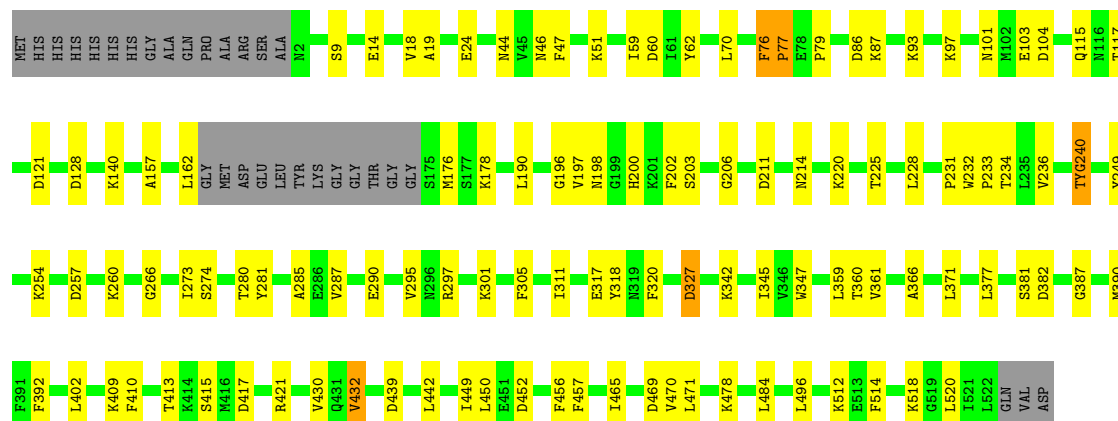
Chain B: 





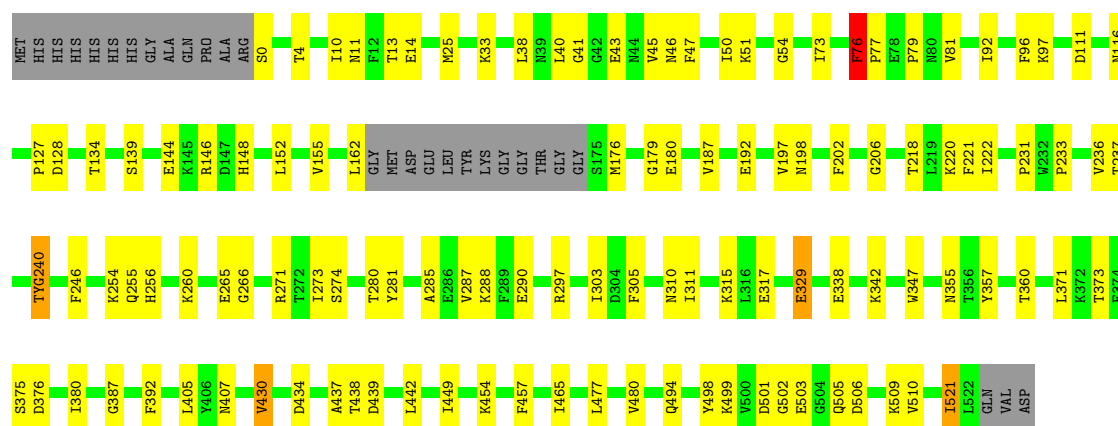
• Molecule 1: iNicSnFR 3.0 Fluorescent Nicotine Sensor

Chain C: 73% 20% • 6%



• Molecule 1: iNicSnFR 3.0 Fluorescent Nicotine Sensor

Chain D: 73% 21% • 5%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	81.19Å 86.86Å 91.83Å 75.72° 85.56° 85.28°	Depositor
Resolution (Å)	40.39 – 2.50 40.39 – 2.50	Depositor EDS
% Data completeness (in resolution range)	98.2 (40.39-2.50) 98.6 (40.39-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.89 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.17.1 _3660	Depositor
R, R_{free}	0.232 , 0.275 0.231 , 0.273	Depositor DCC
R_{free} test set	2007 reflections (2.43%)	wwPDB-VP
Wilson B-factor (Å ²)	40.7	Xtriage
Anisotropy	0.252	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 31.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16649	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.76% 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: CRO

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	0.22	0/4103	0.47	0/5547
1	B	0.22	1/4098 (0.0%)	0.48	0/5540
1	C	0.19	0/4087	0.46	0/5525
1	D	0.21	0/4104	0.48	0/5548
All	All	0.21	1/16392 (0.0%)	0.47	0/22160

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	1
1	C	0	1
1	D	0	1
All	All	0	4

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	322	PRO	CA-C	-5.30	1.49	1.51

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	76	PHE	Peptide

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Mol	Chain	Res	Type	Group
1	B	76	PHE	Peptide
1	C	76	PHE	Peptide
1	D	76	PHE	Peptide

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	4040	0	4022	84	0
1	B	4035	0	4019	92	0
1	C	4024	0	4005	77	0
1	D	4038	0	4024	73	0
2	A	134	0	0	2	0
2	B	112	0	0	6	0
2	C	147	0	0	7	0
2	D	119	0	0	1	0
All	All	16649	0	16070	316	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 10.

All (316) 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:D:76:PHE:HE2	1:D:338:GLU:HG3	1.36	0.91
1:B:361:VAL:HG12	1:B:432:VAL:HG12	1.62	0.81
1:A:18:VAL:HG12	1:A:63:VAL:HG21	1.66	0.78
1:A:442:LEU:HD13	1:A:449:ILE:HD11	1.69	0.74
1:C:442:LEU:HD13	1:C:449:ILE:HD11	1.69	0.73
1:B:76:PHE:HE2	1:B:338:GLU:HG3	1.54	0.71
1:D:4:THR:HG22	1:D:33:LYS:HB2	1.72	0.71
1:B:70:LEU:HG	1:B:335:VAL:HG13	1.72	0.70
1:C:51:LYS:NZ	1:C:104:ASP:O	2.26	0.69
1:D:76:PHE:CE2	1:D:338:GLU:HG3	2.24	0.69
1:C:390:MET:HG2	1:C:417:ASP:OD2	1.94	0.67
1:B:198:ASN:ND2	1:B:305:PHE:O	2.27	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:76:PHE:CE2	1:B:338:GLU:HG3	2.30	0.67
1:C:361:VAL:HG22	1:C:450:LEU:HD21	1.77	0.67
1:B:176:MET:HE1	1:D:40:LEU:HD13	1.77	0.66
1:B:415:SER:OG	2:B:601:HOH:O	2.13	0.66
1:A:466:ILE:HD13	1:A:471:LEU:HD13	1.78	0.66
1:D:192:GLU:OE2	1:D:297:ARG:NH1	2.29	0.65
1:A:235:LEU:O	1:A:239:LEU:HG	1.97	0.65
1:D:14:GLU:OE1	1:D:355:ASN:ND2	2.29	0.65
1:A:496:LEU:HD22	1:A:510:VAL:HG13	1.79	0.64
1:C:198:ASN:ND2	1:C:305:PHE:O	2.30	0.64
1:D:439:ASP:HB3	1:D:442:LEU:HG	1.78	0.64
1:A:2:ASN:ND2	1:D:0:SER:O	2.31	0.64
1:D:25:MET:HA	1:D:521:ILE:HD11	1.81	0.63
1:C:211:ASP:OD2	1:C:214:ASN:ND2	2.25	0.63
1:A:317:GLU:OE1	1:A:342:LYS:NZ	2.31	0.63
1:D:134:THR:HG23	1:D:155:VAL:HG22	1.80	0.62
1:C:387:GLY:N	1:C:430:VAL:HG21	2.15	0.62
1:B:439:ASP:HB3	1:B:442:LEU:HG	1.81	0.61
1:D:198:ASN:ND2	1:D:305:PHE:O	2.33	0.61
1:C:76:PHE:CD2	1:C:77:PRO:HD3	2.35	0.61
1:A:70:LEU:HD22	1:A:76:PHE:CD2	2.36	0.61
1:C:471:LEU:HD21	1:C:478:LYS:HE2	1.83	0.60
1:D:144:GLU:OE1	1:D:220:LYS:NZ	2.34	0.60
1:A:341:ARG:CZ	1:B:390:MET:HE2	2.32	0.59
1:D:499:LYS:HD2	1:D:510:VAL:HG13	1.85	0.59
1:C:290:GLU:OE2	1:C:297:ARG:NH1	2.29	0.59
1:C:317:GLU:OE1	1:C:342:LYS:NZ	2.36	0.59
1:D:176:MET:HE2	1:D:255:GLN:HB3	1.84	0.59
1:D:357:TYR:OH	1:D:434:ASP:OD2	2.17	0.58
1:A:392:PHE:HE1	1:A:457:PHE:HZ	1.51	0.58
1:A:514:PHE:CZ	1:A:518:LYS:HD2	2.38	0.58
1:B:340:LYS:NZ	1:B:344:ASN:OD1	2.32	0.58
1:D:329:GLU:N	1:D:329:GLU:OE1	2.37	0.58
1:B:513:GLU:HA	1:B:516:LYS:HE2	1.86	0.57
1:B:513:GLU:HA	1:B:516:LYS:CE	2.34	0.57
1:B:509:LYS:HD3	1:B:512:LYS:HD3	1.87	0.57
1:B:60:ASP:HB3	1:B:470:VAL:HG21	1.86	0.56
1:C:387:GLY:H	1:C:430:VAL:HG21	1.70	0.56
1:A:70:LEU:CD2	1:A:335:VAL:HG13	2.34	0.56
1:D:509:LYS:HD2	1:D:509:LYS:N	2.21	0.56
1:A:439:ASP:HB3	1:A:442:LEU:HG	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:162:LEU:H	1:D:162:LEU:HD23	1.71	0.56
1:B:477:LEU:HA	1:B:480:VAL:HG22	1.87	0.56
1:B:70:LEU:HD11	1:B:331:ALA:HB1	1.86	0.55
1:B:26:ILE:O	1:B:30:THR:OG1	2.20	0.55
1:A:231:PRO:O	1:A:234:THR:OG1	2.23	0.55
1:C:60:ASP:HB3	1:C:470:VAL:HG21	1.88	0.55
1:A:513:GLU:O	1:A:517:GLU:HG2	2.06	0.55
1:B:514:PHE:CZ	1:B:518:LYS:HD2	2.41	0.55
1:B:498:TYR:CE1	1:B:502:GLY:HA3	2.42	0.55
1:D:233:PRO:HA	1:D:236:VAL:HG23	1.89	0.55
1:B:233:PRO:HA	1:B:236:VAL:HG23	1.89	0.55
1:A:51:LYS:NZ	1:A:104:ASP:O	2.38	0.54
1:C:14:GLU:O	1:C:18:VAL:HG22	2.08	0.54
1:D:392:PHE:HE1	1:D:457:PHE:HZ	1.56	0.54
1:A:471:LEU:HD21	1:A:478:LYS:HE2	1.90	0.54
1:B:347:TRP:HA	1:B:465:ILE:HG22	1.89	0.54
1:B:503:GLU:OE1	1:B:505:GLN:NE2	2.41	0.54
1:A:347:TRP:HA	1:A:465:ILE:HG22	1.90	0.54
1:D:506:ASP:HB3	1:D:509:LYS:HD3	1.89	0.54
1:B:438:THR:HA	1:B:494:GLN:HG3	1.91	0.53
1:A:361:VAL:HG22	1:A:450:LEU:HD21	1.90	0.53
1:A:134:THR:HG23	1:A:155:VAL:HG22	1.91	0.53
1:A:143:ASN:ND2	1:B:501:ASP:O	2.36	0.52
1:B:329:GLU:N	1:B:329:GLU:OE1	2.42	0.52
1:C:274:SER:HA	1:C:280:THR:HG22	1.91	0.52
1:C:162:LEU:HD23	1:C:162:LEU:H	1.74	0.52
1:C:392:PHE:HE1	1:C:457:PHE:HZ	1.57	0.52
1:A:198:ASN:ND2	1:A:305:PHE:O	2.40	0.52
1:B:115:GLN:HB3	2:B:604:HOH:O	2.09	0.52
1:B:273:ILE:HB	1:B:281:TYR:HB2	1.92	0.52
1:A:452:ASP:OD2	1:A:456:PHE:N	2.43	0.52
1:C:70:LEU:HB3	1:C:76:PHE:HB3	1.92	0.52
1:A:122:GLY:O	2:A:601:HOH:O	2.19	0.52
1:B:162:LEU:HD23	1:B:162:LEU:H	1.74	0.52
1:A:228:LEU:HD22	1:A:232:TRP:CE2	2.46	0.51
1:A:139:SER:HB2	1:B:418:MET:HB3	1.91	0.51
1:C:231:PRO:O	1:C:234:THR:OG1	2.25	0.51
1:B:492:GLU:O	1:B:496:LEU:HD13	2.10	0.51
1:C:70:LEU:HD22	1:C:76:PHE:CD2	2.45	0.51
1:B:175:SER:OG	1:B:177:SER:O	2.29	0.51
1:A:492:GLU:OE1	1:A:518:LYS:NZ	2.43	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:231:PRO:HD3	1:D:311:ILE:O	2.11	0.50
1:A:290:GLU:OE2	1:A:297:ARG:NH1	2.36	0.50
1:A:499:LYS:HD3	1:A:510:VAL:HG22	1.93	0.50
1:D:197:VAL:HG23	1:D:202:PHE:HE1	1.76	0.50
1:D:407:ASN:O	1:D:407:ASN:ND2	2.44	0.50
1:C:228:LEU:HD22	1:C:232:TRP:CE2	2.47	0.50
1:D:148:HIS:HB3	1:D:222:ILE:HD13	1.93	0.50
1:D:273:ILE:HB	1:D:281:TYR:HB2	1.92	0.50
1:A:211:ASP:OD2	1:A:214:ASN:ND2	2.39	0.50
1:D:477:LEU:HA	1:D:480:VAL:HG22	1.93	0.50
1:B:197:VAL:HG23	1:B:202:PHE:HE1	1.77	0.49
1:D:371:LEU:HD11	1:D:380:ILE:HD11	1.93	0.49
1:C:233:PRO:HA	1:C:236:VAL:HG23	1.94	0.49
1:D:266:GLY:HA3	1:D:287:VAL:O	2.12	0.49
1:D:438:THR:HA	1:D:494:GLN:HG3	1.94	0.49
1:D:97:LYS:HG2	1:D:111:ASP:OD1	2.13	0.49
1:A:76:PHE:CG	1:A:77:PRO:N	2.80	0.49
1:B:157:ALA:HB3	1:B:249:TYR:HA	1.94	0.49
1:B:507:PRO:HA	1:B:510:VAL:HG12	1.94	0.49
1:C:347:TRP:HA	1:C:465:ILE:HG22	1.93	0.49
1:A:57:ASN:ND2	1:C:128:ASP:OD1	2.43	0.49
1:C:439:ASP:HB3	1:C:442:LEU:HG	1.92	0.49
1:D:236:VAL:O	1:D:240:CRO:N1	2.45	0.49
1:A:438:THR:HA	1:A:494:GLN:HG3	1.94	0.49
1:B:390:MET:HG2	1:B:417:ASP:OD1	2.13	0.49
1:A:240:CRO:HA1	1:A:243:VAL:HG22	1.95	0.49
1:A:361:VAL:HG12	1:A:432:VAL:HG22	1.94	0.48
1:C:327:ASP:OD2	1:C:327:ASP:N	2.38	0.48
1:A:221:PHE:CZ	1:A:239:LEU:HD13	2.49	0.48
1:A:495:LYS:HE3	1:A:499:LYS:HE2	1.94	0.48
1:A:59:ILE:HD12	1:A:62:TYR:HB3	1.95	0.48
1:A:290:GLU:CD	1:A:297:ARG:HH12	2.20	0.48
1:B:403:GLN:HG3	1:B:408:PHE:HB2	1.95	0.48
1:C:197:VAL:HG23	1:C:202:PHE:HE1	1.79	0.48
1:C:285:ALA:HA	1:C:297:ARG:O	2.13	0.48
1:D:79:PRO:HG3	1:D:97:LYS:O	2.13	0.48
1:B:101:ASN:OD1	2:B:602:HOH:O	2.20	0.48
1:B:357:TYR:OH	1:B:434:ASP:OD2	2.21	0.48
1:D:405:LEU:HG	1:D:454:LYS:HE2	1.95	0.48
1:B:74:LEU:HD13	1:B:76:PHE:CE2	2.48	0.48
1:B:414:LYS:HD3	1:B:416:MET:HE3	1.94	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:93:LYS:HG2	1:B:115:GLN:HG2	1.96	0.48
1:B:384:LEU:HB3	1:B:431:GLN:HB2	1.95	0.48
1:D:303:ILE:HD11	2:D:688:HOH:O	2.14	0.48
1:C:59:ILE:HD12	1:C:62:TYR:HB3	1.96	0.47
1:A:299:GLU:OE2	1:A:301:LYS:HG2	2.14	0.47
1:B:240:CRO:HD1	1:B:240:CRO:N2	2.29	0.47
1:D:41:GLY:O	1:D:45:VAL:HB	2.13	0.47
1:D:13:THR:HG21	1:D:501:ASP:OD1	2.15	0.47
1:A:240:CRO:N2	1:A:240:CRO:HD1	2.29	0.47
1:D:347:TRP:HA	1:D:465:ILE:HG22	1.96	0.47
1:A:206:GLY:HA2	1:A:220:LYS:O	2.15	0.47
1:B:351:LEU:HB3	1:B:488:ILE:HD12	1.97	0.47
1:B:231:PRO:HD3	1:B:311:ILE:O	2.15	0.47
1:C:231:PRO:HD3	1:C:311:ILE:O	2.14	0.47
1:A:60:ASP:HB3	1:A:470:VAL:HG21	1.97	0.47
1:A:380:ILE:HG12	1:A:380:ILE:O	2.14	0.47
1:B:392:PHE:HE1	1:B:457:PHE:HZ	1.63	0.47
1:A:231:PRO:HD3	1:A:311:ILE:O	2.15	0.47
1:B:285:ALA:HA	1:B:297:ARG:O	2.15	0.46
1:C:240:CRO:N2	1:C:240:CRO:HD1	2.30	0.46
1:B:231:PRO:O	1:B:234:THR:OG1	2.31	0.46
1:C:381:SER:O	1:C:410:PHE:HA	2.16	0.46
1:C:514:PHE:CZ	1:C:518:LYS:HD2	2.50	0.46
1:D:25:MET:SD	1:D:521:ILE:HD13	2.56	0.46
1:A:62:TYR:CZ	1:A:465:ILE:HD11	2.51	0.46
1:B:134:THR:HG23	1:B:155:VAL:HG22	1.97	0.46
1:C:266:GLY:HA3	1:C:287:VAL:O	2.16	0.46
1:C:382:ASP:HA	1:C:409:LYS:O	2.15	0.46
1:C:115:GLN:HB3	2:C:601:HOH:O	2.16	0.46
1:A:197:VAL:HG23	1:A:202:PHE:HE1	1.81	0.46
1:A:221:PHE:CE2	1:A:239:LEU:HD13	2.50	0.46
1:A:273:ILE:HB	1:A:281:TYR:HB2	1.97	0.46
1:C:178:LYS:HE2	2:C:717:HOH:O	2.16	0.46
1:B:20:ASN:O	1:B:24:GLU:HG2	2.16	0.46
1:B:198:ASN:N	2:B:608:HOH:O	2.49	0.46
1:B:417:ASP:OD2	2:B:603:HOH:O	2.21	0.46
1:D:11:ASN:HA	1:D:38:LEU:HD11	1.98	0.46
1:D:47:PHE:CE2	1:D:51:LYS:HD2	2.51	0.46
1:A:484:LEU:HD23	1:A:520:LEU:HD13	1.98	0.45
1:B:97:LYS:HE2	1:B:111:ASP:OD1	2.17	0.45
1:B:337:LYS:HD2	1:B:341:ARG:NH2	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:179:GLY:HA3	1:D:260:LYS:O	2.16	0.45
1:D:206:GLY:HA3	1:D:221:PHE:CD1	2.52	0.45
1:B:5:VAL:HG23	1:B:34:VAL:HG22	1.98	0.45
1:C:257:ASP:OD2	1:C:260:LYS:NZ	2.49	0.45
1:A:70:LEU:HD11	1:A:331:ALA:HB1	1.98	0.45
1:A:179:GLY:HA3	1:A:260:LYS:O	2.15	0.45
1:A:516:LYS:HG3	1:A:521:ILE:HD11	1.99	0.45
1:B:512:LYS:O	1:B:516:LYS:HG3	2.17	0.45
1:D:92:ILE:HG12	1:D:116:ASN:HB2	1.97	0.45
1:A:62:TYR:CE1	1:A:465:ILE:HD11	2.52	0.45
1:B:79:PRO:HG3	1:B:97:LYS:O	2.17	0.45
1:C:366:ALA:HA	1:C:371:LEU:HD12	1.99	0.45
1:A:270:GLU:HG2	1:A:284:ARG:HG3	1.98	0.45
1:B:148:HIS:HB3	1:B:222:ILE:HD13	1.99	0.45
1:C:115:GLN:NE2	2:C:602:HOH:O	2.49	0.45
1:D:274:SER:HA	1:D:280:THR:HG22	1.99	0.45
1:D:373:THR:HG23	1:D:375:SER:H	1.82	0.45
1:B:32:LEU:HD11	1:B:474:HIS:CG	2.51	0.45
1:C:190:LEU:O	1:C:295:VAL:HA	2.17	0.45
1:A:233:PRO:HA	1:A:236:VAL:HG23	1.99	0.44
1:B:393:LEU:HA	1:B:399:TYR:HB3	1.99	0.44
1:A:18:VAL:CG1	1:A:63:VAL:HG21	2.43	0.44
1:C:496:LEU:HD11	1:C:514:PHE:CD1	2.52	0.44
1:A:135:GLN:HG2	1:A:323:PRO:HB3	2.00	0.44
1:A:415:SER:OG	2:A:602:HOH:O	2.20	0.44
1:A:3:ASP:OD1	1:B:88:GLN:HG3	2.17	0.44
1:C:178:LYS:HG2	2:C:683:HOH:O	2.16	0.44
1:B:117:THR:OG1	2:B:604:HOH:O	2.21	0.44
1:B:331:ALA:O	1:B:335:VAL:HG22	2.18	0.44
1:D:92:ILE:CG1	1:D:116:ASN:HB2	2.47	0.44
1:D:265:GLU:CD	1:D:265:GLU:H	2.26	0.44
1:D:285:ALA:HA	1:D:297:ARG:O	2.18	0.44
1:B:322:PRO:HA	1:B:323:PRO:HD3	1.88	0.44
1:C:377:LEU:HD21	1:C:432:VAL:HG11	2.00	0.44
1:C:79:PRO:HG3	1:C:97:LYS:O	2.18	0.44
1:D:81:VAL:HG13	1:D:96:PHE:CD1	2.53	0.44
1:A:181:GLU:O	1:A:184:THR:HG23	2.18	0.44
1:A:381:SER:O	1:A:410:PHE:HA	2.17	0.44
1:D:357:TYR:CE1	1:D:434:ASP:HB2	2.53	0.44
1:D:498:TYR:CE1	1:D:502:GLY:HA3	2.52	0.44
1:C:196:GLY:HA2	1:C:200:HIS:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:345:ILE:HG21	1:C:465:ILE:HD12	2.00	0.43
1:C:361:VAL:HG12	1:C:432:VAL:HG13	2.00	0.43
1:A:496:LEU:HD23	1:A:499:LYS:HD2	1.99	0.43
1:C:273:ILE:HB	1:C:281:TYR:HB2	1.99	0.43
1:C:140:LYS:HE2	1:C:318:TYR:CZ	2.53	0.43
1:C:421:ARG:NH1	2:C:608:HOH:O	2.34	0.43
1:D:128:ASP:H	1:D:256:HIS:CD2	2.35	0.43
1:D:310:ASN:HA	1:D:315:LYS:HB2	2.00	0.43
1:B:50:ILE:HA	1:B:59:ILE:O	2.19	0.43
1:D:187:VAL:HB	1:D:246:PHE:HE1	1.83	0.43
1:C:360:THR:HG22	1:C:449:ILE:HA	2.01	0.43
1:C:469:ASP:OD1	1:C:470:VAL:N	2.50	0.43
1:B:76:PHE:CD1	1:B:76:PHE:N	2.86	0.43
1:B:274:SER:HA	1:B:280:THR:HG22	2.01	0.43
1:B:357:TYR:CE1	1:B:434:ASP:HB2	2.54	0.43
1:B:495:LYS:HE2	1:B:495:LYS:HB3	1.77	0.43
1:B:13:THR:HG21	1:B:501:ASP:OD2	2.19	0.43
1:C:452:ASP:OD2	1:C:457:PHE:N	2.34	0.43
1:A:35:VAL:HG21	1:C:87:LYS:O	2.19	0.43
1:B:177:SER:OG	1:B:261:SER:HA	2.19	0.43
1:C:360:THR:HG22	1:C:449:ILE:HG13	2.01	0.43
1:B:2:ASN:OD1	1:B:31:ASP:HB3	2.19	0.42
1:C:70:LEU:HD22	1:C:76:PHE:CG	2.54	0.42
1:C:176:MET:HG2	1:C:254:LYS:HD3	2.01	0.42
1:D:10:ILE:HG23	1:D:46:ASN:ND2	2.33	0.42
1:A:266:GLY:HA3	1:A:287:VAL:O	2.19	0.42
1:C:301:LYS:HA	1:C:301:LYS:HD3	1.82	0.42
1:B:47:PHE:O	1:B:50:ILE:HG22	2.19	0.42
1:D:240:CRO:HD1	1:D:240:CRO:N2	2.34	0.42
1:B:426:ASP:OD2	1:B:445:HIS:ND1	2.44	0.42
1:B:62:TYR:CZ	1:B:465:ILE:HD11	2.55	0.42
1:D:180:GLU:OE2	1:D:254:LYS:HD3	2.19	0.42
1:A:79:PRO:HG3	1:A:97:LYS:O	2.20	0.42
1:B:187:VAL:HB	1:B:246:PHE:HE1	1.84	0.42
1:C:46:ASN:ND2	2:C:631:HOH:O	2.52	0.42
1:D:317:GLU:OE2	1:D:342:LYS:NZ	2.47	0.42
1:A:187:VAL:HB	1:A:246:PHE:HE1	1.84	0.42
1:B:16:ILE:HG12	1:B:38:LEU:HD11	2.01	0.42
1:C:86:ASP:HB2	1:C:93:LYS:HG3	2.01	0.42
1:C:103:GLU:HG3	1:C:317:GLU:HG3	2.02	0.42
1:A:153:GLU:HB2	1:A:217:LEU:HB2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:354:ASN:HB2	1:B:461:TYR:CE1	2.55	0.42
1:C:117:THR:OG1	2:C:601:HOH:O	2.21	0.42
1:A:366:ALA:HA	1:A:371:LEU:HD12	2.02	0.42
1:B:87:LYS:NZ	1:D:54:GLY:O	2.49	0.42
1:C:44:ASN:HB2	1:C:47:PHE:HB3	2.01	0.42
1:C:452:ASP:OD2	1:C:456:PHE:N	2.53	0.42
1:D:146:ARG:HB2	1:D:222:ILE:HD12	2.01	0.42
1:A:36:ARG:HD2	1:C:121:ASP:O	2.20	0.41
1:B:162:LEU:HB3	1:B:252:HIS:CG	2.54	0.41
1:B:390:MET:HE1	1:B:394:GLU:CD	2.45	0.41
1:D:237:THR:O	1:D:271:ARG:NH1	2.46	0.41
1:A:187:VAL:HB	1:A:246:PHE:CE1	2.55	0.41
1:A:206:GLY:HA3	1:A:221:PHE:CD1	2.55	0.41
1:A:360:THR:HG22	1:A:449:ILE:HA	2.02	0.41
1:C:157:ALA:HB3	1:C:249:TYR:HA	2.01	0.41
1:C:203:SER:HB2	1:C:225:THR:HG23	2.02	0.41
1:D:47:PHE:O	1:D:50:ILE:HG22	2.21	0.41
1:D:437:ALA:O	1:D:494:GLN:NE2	2.39	0.41
1:A:385:ILE:HG22	1:A:430:VAL:HG22	2.02	0.41
1:A:423:THR:HG22	1:A:427:ASN:ND2	2.35	0.41
1:B:126:LEU:HD23	1:B:126:LEU:HA	1.81	0.41
1:B:187:VAL:HB	1:B:246:PHE:CE1	2.55	0.41
1:B:9:SER:O	1:B:38:LEU:HA	2.19	0.41
1:C:206:GLY:HA2	1:C:220:LYS:O	2.20	0.41
1:C:390:MET:SD	1:C:415:SER:OG	2.70	0.41
1:D:127:PRO:HB3	1:D:256:HIS:O	2.21	0.41
1:A:18:VAL:HG11	1:A:63:VAL:HG11	2.01	0.41
1:B:513:GLU:HA	1:B:516:LYS:HE3	2.01	0.41
1:C:228:LEU:HD22	1:C:232:TRP:CD2	2.56	0.41
1:A:61:ILE:HG12	1:A:466:ILE:HB	2.03	0.41
1:B:111:ASP:O	1:B:273:ILE:HA	2.20	0.41
1:B:509:LYS:O	1:B:512:LYS:HG2	2.21	0.41
1:D:198:ASN:OD1	1:D:305:PHE:HB2	2.21	0.41
1:D:387:GLY:HA3	1:D:430:VAL:HG21	2.02	0.41
1:B:337:LYS:HB2	1:B:337:LYS:HE2	1.82	0.41
1:B:458:PRO:HG2	1:B:460:TYR:HE2	1.85	0.41
1:B:498:TYR:CD1	1:B:502:GLY:HA3	2.56	0.41
1:C:24:GLU:HG3	1:C:512:LYS:HG2	2.03	0.41
1:C:359:LEU:HD11	1:C:402:LEU:HD21	2.03	0.41
1:D:288:LYS:NZ	1:D:290:GLU:OE2	2.44	0.41
1:D:503:GLU:OE1	1:D:505:GLN:NE2	2.52	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:9:SER:HB2	1:A:19:ALA:CB	2.51	0.41
1:A:257:ASP:OD2	1:A:260:LYS:NZ	2.43	0.41
1:A:358:THR:OG1	1:A:449:ILE:HG23	2.21	0.41
1:A:499:LYS:HB3	1:A:510:VAL:HG21	2.02	0.41
1:A:70:LEU:HD23	1:A:335:VAL:HG13	2.02	0.40
1:A:86:ASP:HB2	1:A:93:LYS:HG3	2.03	0.40
1:C:361:VAL:CG1	1:C:432:VAL:HG13	2.51	0.40
1:C:512:LYS:HB3	1:C:512:LYS:HE2	1.83	0.40
1:D:152:LEU:HD12	1:D:218:THR:HG22	2.04	0.40
1:D:360:THR:HG22	1:D:449:ILE:HA	2.03	0.40
1:C:9:SER:HB2	1:C:19:ALA:CB	2.51	0.40
1:C:484:LEU:HD23	1:C:520:LEU:HD13	2.03	0.40
1:D:187:VAL:HB	1:D:246:PHE:CE1	2.57	0.40
1:D:373:THR:HG22	1:D:376:ASP:CG	2.45	0.40
1:A:406:TYR:OH	1:A:452:ASP:OD1	2.36	0.40
1:B:48:GLU:HA	1:B:51:LYS:NZ	2.36	0.40
1:B:280:THR:O	1:B:302:GLY:HA2	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	502/537 (94%)	492 (98%)	9 (2%)	1 (0%)	43	63
1	B	502/537 (94%)	493 (98%)	8 (2%)	1 (0%)	43	63
1	C	500/537 (93%)	491 (98%)	8 (2%)	1 (0%)	43	63
1	D	503/537 (94%)	492 (98%)	9 (2%)	2 (0%)	30	49
All	All	2007/2148 (93%)	1968 (98%)	34 (2%)	5 (0%)	43	63

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	77	PRO
1	D	77	PRO
1	C	77	PRO
1	A	76	PHE
1	D	76	PHE

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	440/459 (96%)	431 (98%)	9 (2%)	48	75
1	B	439/459 (96%)	426 (97%)	13 (3%)	36	64
1	C	438/459 (95%)	433 (99%)	5 (1%)	65	84
1	D	440/459 (96%)	433 (98%)	7 (2%)	55	79
All	All	1757/1836 (96%)	1723 (98%)	34 (2%)	50	76

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	43	GLU
1	A	44	ASN
1	A	95	ASN
1	A	152	LEU
1	A	320	PHE
1	A	326	THR
1	A	335	VAL
1	A	510	VAL
1	A	516	LYS
1	B	35	VAL
1	B	38	LEU
1	B	103	GLU
1	B	117	THR
1	B	156	THR
1	B	299	GLU
1	B	326	THR
1	B	335	VAL

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Mol	Chain	Res	Type
1	B	368	GLN
1	B	380	ILE
1	B	412	HIS
1	B	430	VAL
1	B	509	LYS
1	C	101	ASN
1	C	320	PHE
1	C	327	ASP
1	C	413	THR
1	C	432	VAL
1	D	43	GLU
1	D	73	ILE
1	D	139[A]	SER
1	D	139[B]	SER
1	D	329	GLU
1	D	430	VAL
1	D	521	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	11	ASN
1	A	95	ASN
1	A	403	GLN
1	B	44	ASN
1	B	68	HIS
1	B	412	HIS
1	B	505	GLN
1	C	101	ASN
1	C	115	GLN
1	C	319	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

4 non-standard protein/DNA/RNA residues 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
1	CRO	D	240	1	22,23,24	2.79	7 (31%)	30,32,34	2.23	6 (20%)
1	CRO	A	240	1	22,23,24	2.73	7 (31%)	30,32,34	2.28	9 (30%)
1	CRO	C	240	1	22,23,24	2.77	7 (31%)	30,32,34	2.23	8 (26%)
1	CRO	B	240	1	22,23,24	2.78	7 (31%)	30,32,34	2.31	8 (26%)

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	CRO	D	240	1	-	3/12/31/32	0/2/2/2
1	CRO	A	240	1	-	3/12/31/32	0/2/2/2
1	CRO	C	240	1	-	6/12/31/32	0/2/2/2
1	CRO	B	240	1	-	2/12/31/32	0/2/2/2

All (28) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	240	CRO	C1-N2	7.54	1.43	1.32
1	C	240	CRO	C1-N2	7.49	1.43	1.32
1	B	240	CRO	C1-N2	7.44	1.43	1.32
1	A	240	CRO	C1-N2	7.38	1.43	1.32
1	B	240	CRO	CA2-C2	5.68	1.54	1.48
1	C	240	CRO	CA2-C2	5.65	1.54	1.48
1	D	240	CRO	CA2-C2	5.63	1.54	1.48
1	A	240	CRO	CA2-C2	5.57	1.54	1.48
1	D	240	CRO	C1-N3	4.89	1.45	1.37
1	C	240	CRO	C1-N3	4.83	1.45	1.37
1	B	240	CRO	C1-N3	4.79	1.45	1.37
1	B	240	CRO	C2-N3	4.58	1.50	1.40
1	A	240	CRO	C1-N3	4.56	1.44	1.37
1	D	240	CRO	C2-N3	4.54	1.50	1.40
1	A	240	CRO	C2-N3	4.41	1.50	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	240	CRO	C2-N3	4.23	1.49	1.40
1	A	240	CRO	CA2-N2	2.67	1.44	1.38
1	B	240	CRO	CA2-N2	2.66	1.44	1.38
1	D	240	CRO	CA2-N2	2.62	1.44	1.38
1	C	240	CRO	CA2-N2	2.60	1.44	1.38
1	D	240	CRO	CA3-N3	2.38	1.51	1.47
1	B	240	CRO	CA1-N1	-2.37	1.40	1.47
1	C	240	CRO	CA3-N3	2.37	1.51	1.47
1	B	240	CRO	CA3-N3	2.34	1.51	1.47
1	A	240	CRO	CA3-N3	2.33	1.51	1.47
1	C	240	CRO	CA1-N1	-2.32	1.41	1.47
1	D	240	CRO	CA1-N1	-2.28	1.41	1.47
1	A	240	CRO	CA1-N1	-2.17	1.41	1.47

All (31) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	240	CRO	O2-C2-CA2	-6.95	126.58	131.02
1	D	240	CRO	O2-C2-CA2	-6.89	126.63	131.02
1	A	240	CRO	O2-C2-CA2	-6.33	126.98	131.02
1	C	240	CRO	O2-C2-CA2	-5.98	127.20	131.02
1	D	240	CRO	C2-N3-C1	-5.69	105.43	108.07
1	C	240	CRO	C2-N3-C1	-5.60	105.47	108.07
1	B	240	CRO	C2-N3-C1	-5.56	105.49	108.07
1	A	240	CRO	C2-N3-C1	-5.01	105.75	108.07
1	A	240	CRO	CA2-N2-C1	4.37	109.22	105.80
1	B	240	CRO	CA2-N2-C1	4.30	109.16	105.80
1	C	240	CRO	CA2-C2-N3	4.13	106.97	103.50
1	D	240	CRO	CA2-N2-C1	4.03	108.95	105.80
1	B	240	CRO	CA2-C2-N3	3.96	106.83	103.50
1	D	240	CRO	CA2-C2-N3	3.94	106.81	103.50
1	C	240	CRO	CA2-N2-C1	3.93	108.87	105.80
1	A	240	CRO	CA2-C2-N3	3.91	106.78	103.50
1	A	240	CRO	C2-CA2-N2	-3.66	106.33	108.95
1	B	240	CRO	C2-CA2-N2	-3.50	106.44	108.95
1	B	240	CRO	CG2-CB2-CA2	-3.41	125.81	129.87
1	C	240	CRO	C2-CA2-N2	-3.39	106.52	108.95
1	C	240	CRO	CG2-CB2-CA2	-3.36	125.87	129.87
1	A	240	CRO	CG2-CB2-CA2	-3.22	126.04	129.87
1	D	240	CRO	C2-CA2-N2	-3.21	106.65	108.95
1	D	240	CRO	CG2-CB2-CA2	-2.39	127.02	129.87
1	A	240	CRO	CA1-C1-N3	-2.39	121.86	124.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	240	CRO	CA3-N3-C2	2.34	128.82	123.67
1	C	240	CRO	CA1-C1-N3	-2.30	121.97	124.69
1	A	240	CRO	CB2-CA2-C2	2.26	125.09	122.36
1	B	240	CRO	CA3-N3-C2	2.18	128.45	123.67
1	B	240	CRO	CB2-CA2-C2	2.08	124.88	122.36
1	C	240	CRO	CB2-CA2-C2	2.02	124.81	122.36

There are no chirality outliers.

All (14) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	240	CRO	C3-CA3-N3-C2
1	B	240	CRO	C3-CA3-N3-C2
1	C	240	CRO	C3-CA3-N3-C2
1	D	240	CRO	C3-CA3-N3-C2
1	C	240	CRO	C2-CA2-CB2-CG2
1	C	240	CRO	N2-CA2-CB2-CG2
1	C	240	CRO	C3-CA3-N3-C1
1	A	240	CRO	C3-CA3-N3-C1
1	B	240	CRO	C3-CA3-N3-C1
1	D	240	CRO	C3-CA3-N3-C1
1	C	240	CRO	N2-C1-CA1-CB1
1	C	240	CRO	N3-C1-CA1-CB1
1	D	240	CRO	N2-C1-CA1-CB1
1	A	240	CRO	N2-CA2-CB2-CG2

There are no ring outliers.

4 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	D	240	CRO	2	0
1	A	240	CRO	2	0
1	C	240	CRO	1	0
1	B	240	CRO	1	0

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	508/537 (94%)	-1.44	0 100 100	20, 36, 55, 80	0
1	B	508/537 (94%)	-1.37	0 100 100	22, 40, 63, 76	0
1	C	506/537 (94%)	-1.42	0 100 100	23, 36, 56, 71	0
1	D	508/537 (94%)	-1.38	0 100 100	19, 41, 63, 103	1 (0%)
All	All	2030/2148 (94%)	-1.40	0 100 100	19, 38, 60, 103	1 (0%)

There are no RSRZ outliers to report.

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. 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(Å ²)	Q<0.9
1	CRO	A	240	22/23	0.99	0.04	19,26,32,34	0
1	CRO	B	240	22/23	0.99	0.04	23,31,36,41	0
1	CRO	C	240	22/23	0.99	0.04	22,27,34,36	0
1	CRO	D	240	22/23	0.99	0.04	31,35,40,42	0

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

6.4 Ligands [i](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.