



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 14, 2017 – 06:51 pm GMT

PDB ID : 4TRY
Title : Structure of BACE1 complex with a HEA-type inhibitor
Authors : Akaji, K.; Teruya, K.; Akiyama, T.; Sanjho, A.; Yamashita, E.; Nakagawa, A.
Deposited on : 2014-06-18
Resolution : 2.75 Å(reported)

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

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

A user guide is available at

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix) : 1.9-1692
EDS : trunk28620
Percentile statistics : 20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : recalc28949

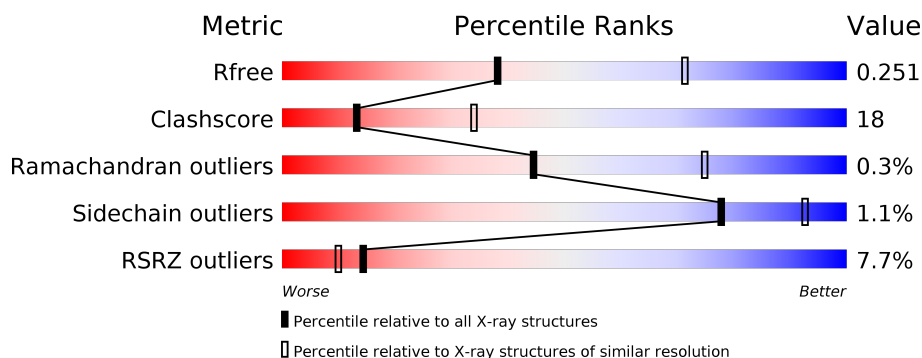
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.75 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}	100719	3666 (2.80-2.72)
Clashscore	112137	4174 (2.80-2.72)
Ramachandran outliers	110173	4103 (2.80-2.72)
Sidechain outliers	110143	4106 (2.80-2.72)
RSRZ outliers	101464	3697 (2.80-2.72)

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	388	<div> <div>7%</div> <div> <div></div> <div>73%</div> <div>26%</div> </div> </div>
1	B	388	<div> <div>7%</div> <div> <div></div> <div>73%</div> <div>27%</div> </div> </div>
1	C	388	<div> <div>10%</div> <div> <div></div> <div>70%</div> <div>28%</div> </div> </div>
2	D	4	<div> <div></div> <div> <div>50%</div> <div>50%</div> </div> </div>
2	E	4	<div> <div></div> <div>100%</div> </div>
2	F	4	<div> <div></div> <div> <div>75%</div> <div>25%</div> </div> </div>

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 9326 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 Beta-secretase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	387	Total	C	N	O	S	0	0	0
			3042	1948	505	575	14			
1	B	388	Total	C	N	O	S	0	0	0
			3050	1954	506	576	14			
1	C	388	Total	C	N	O	S	0	0	0
			3050	1954	506	576	14			

- Molecule 2 is a protein called GLU-ILE-TIH-THC-NVA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	4	Total	C	N	O	S	0	0	0
			47	31	5	9	2			
2	E	4	Total	C	N	O	S	0	0	0
			47	31	5	9	2			
2	F	4	Total	C	N	O	S	0	0	0
			47	31	5	9	2			

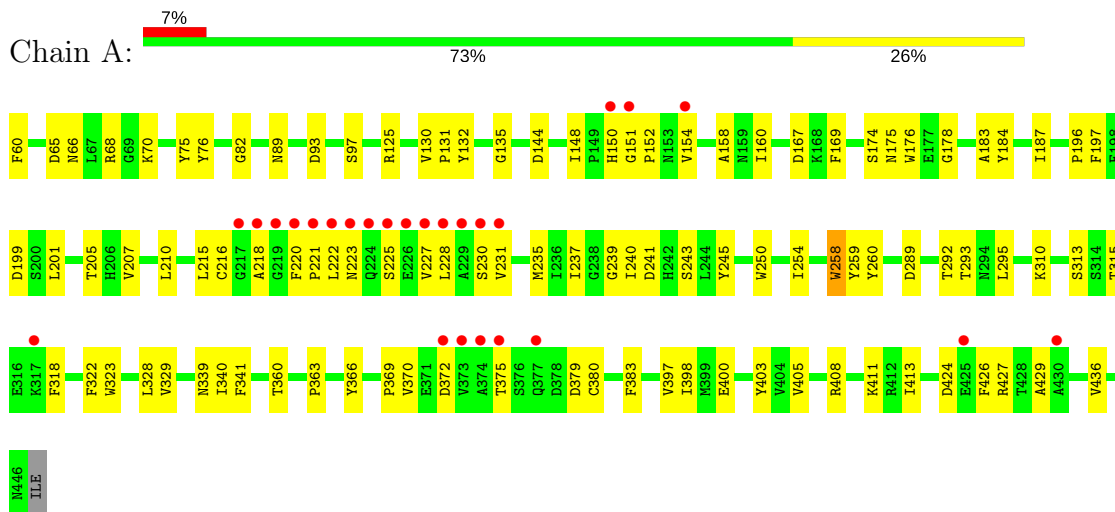
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	10	Total	O	0	0
			10	10		
3	B	15	Total	O	0	0
			15	15		
3	C	16	Total	O	0	0
			16	16		
3	E	1	Total	O	0	0
			1	1		
3	F	1	Total	O	0	0
			1	1		

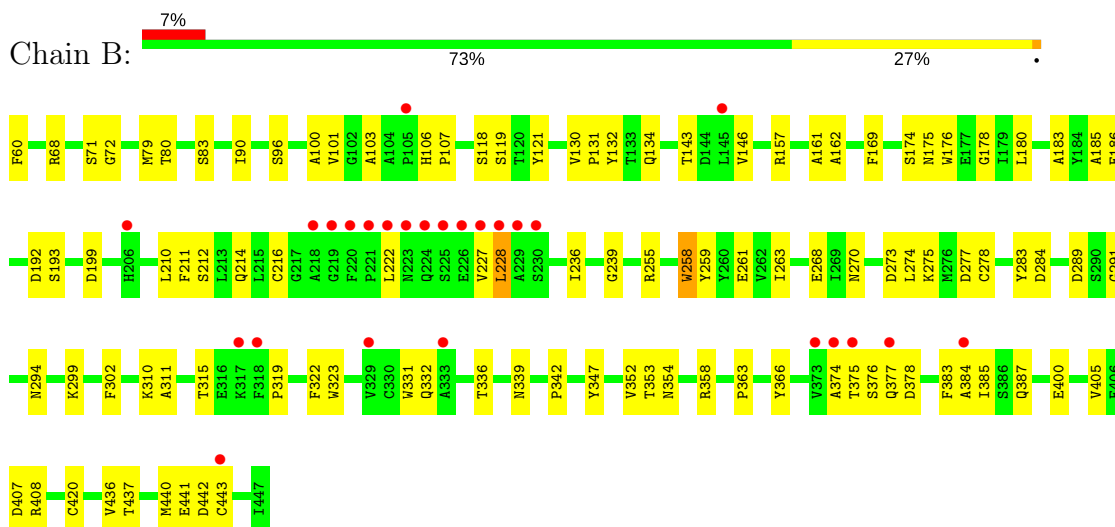
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of 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: Beta-secretase 1

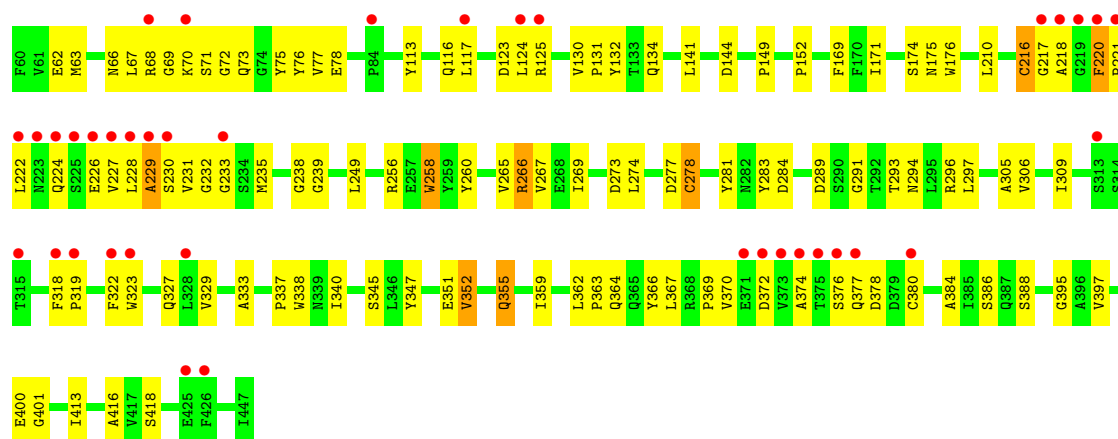


• Molecule 1: Beta-secretase 1



• Molecule 1: Beta-secretase 1

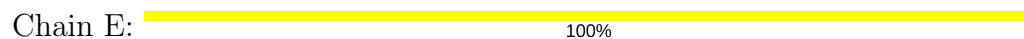




- Molecule 2: GLU-ILE-TIH-THC-NVA



- Molecule 2: GLU-ILE-TIH-THC-NVA



- Molecule 2: GLU-ILE-TIH-THC-NVA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.33Å 102.62Å 101.29Å 90.00° 103.53° 90.00°	Depositor
Resolution (Å)	49.18 – 2.75 49.18 – 2.75	Depositor EDS
% Data completeness (in resolution range)	99.7 (49.18-2.75) 99.7 (49.18-2.75)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.31 (at 2.77Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.179 , 0.252 0.204 , 0.251	Depositor DCC
R_{free} test set	2136 reflections (5.04%)	DCC
Wilson B-factor (Å ²)	57.6	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 37.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	9326	wwPDB-VP
Average B, all atoms (Å ²)	64.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.68% 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: 36D, TIH

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	0.60	0/3120	0.82	1/4243 (0.0%)
1	B	0.63	0/3128	0.81	1/4254 (0.0%)
1	C	0.62	0/3128	0.79	3/4254 (0.1%)
2	D	0.78	0/16	0.46	0/20
2	E	0.18	0/16	0.38	0/20
2	F	0.18	0/16	0.38	0/20
All	All	0.62	0/9424	0.80	5/12811 (0.0%)

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	130	VAL	C-N-CD	5.29	139.52	128.40
1	A	130	VAL	C-N-CD	5.24	139.41	128.40
1	B	130	VAL	C-N-CD	5.21	139.35	128.40
1	C	220	PHE	C-N-CD	5.20	139.32	128.40
1	C	229	ALA	CB-CA-C	5.00	117.61	110.10

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	3042	0	2951	105	0
1	B	3050	0	2964	99	0
1	C	3050	0	2962	135	0
2	D	47	0	43	7	0
2	E	47	0	43	13	0
2	F	47	0	42	8	0
3	A	10	0	0	0	0
3	B	15	0	0	0	0
3	C	16	0	0	1	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
All	All	9326	0	9005	331	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 18.

All (331) 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:C:69:GLY:O	1:C:231:VAL:HB	1.32	1.28
1:C:229:ALA:CA	1:C:230:SER:OG	1.81	1.27
1:C:229:ALA:HA	1:C:230:SER:CB	1.57	1.24
1:B:134:GLN:OE1	2:E:3:TIH:HD	1.40	1.20
1:A:70:LYS:NZ	1:A:222:LEU:HD13	1.65	1.12
1:C:229:ALA:HA	1:C:230:SER:OG	0.92	1.09
1:C:217:GLY:HA2	1:C:231:VAL:HG13	1.32	1.08
1:C:67:LEU:HD11	1:C:77:VAL:HB	1.26	1.07
1:A:235:MET:CE	1:A:237:ILE:HD11	1.88	1.03
1:C:67:LEU:HB2	1:C:233:GLY:O	1.59	1.01
1:A:70:LYS:HZ2	1:A:222:LEU:HD13	1.23	0.98
1:C:217:GLY:HA2	1:C:231:VAL:CG1	1.95	0.96
1:C:210:LEU:HD11	1:C:239:GLY:HA2	1.48	0.96
1:B:310:LYS:HE3	1:B:323:TRP:CD1	2.00	0.95
1:A:65:ASP:OD1	1:A:231:VAL:HG11	1.66	0.95
1:A:235:MET:HE2	1:A:237:ILE:HD11	1.48	0.94
1:C:229:ALA:HB1	1:C:231:VAL:N	1.83	0.92
1:C:210:LEU:HD11	1:C:239:GLY:CA	1.99	0.92
1:C:76:TYR:HE1	1:C:224:GLN:NE2	1.69	0.91
1:C:66:ASN:HD21	1:C:235:MET:H	1.11	0.91
1:A:125:ARG:HH11	1:A:125:ARG:HG3	1.34	0.90
1:C:76:TYR:HE1	1:C:224:GLN:HE21	0.99	0.90
1:B:212:SER:OG	1:B:236:ILE:HB	1.73	0.87

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:216:CYS:O	1:C:231:VAL:HG12	1.75	0.87
1:B:169:PHE:CD1	2:E:4:36D:H9	2.11	0.85
1:C:229:ALA:HB1	1:C:231:VAL:H	1.40	0.85
1:B:270:ASN:ND2	1:B:342:PRO:HB3	1.92	0.84
1:C:305:ALA:O	1:C:309:ILE:HG13	1.78	0.84
1:B:294:ASN:HB3	1:B:384:ALA:O	1.78	0.83
1:C:218:ALA:CB	1:C:220:PHE:O	2.26	0.83
1:A:97:SER:OG	1:A:187:ILE:HD11	1.79	0.83
1:C:222:LEU:HB2	1:C:229:ALA:O	1.78	0.82
1:C:221:PRO:C	1:C:222:LEU:HD23	2.00	0.82
1:C:67:LEU:CB	1:C:233:GLY:O	2.28	0.81
1:C:218:ALA:HB3	1:C:220:PHE:O	1.80	0.81
1:C:66:ASN:ND2	1:C:235:MET:H	1.77	0.80
1:A:426:PHE:CD1	1:B:311:ALA:HB1	2.16	0.80
1:C:351:GLU:H	1:C:355:GLN:HE21	1.30	0.80
1:C:210:LEU:HD11	1:C:239:GLY:N	1.97	0.80
1:A:65:ASP:OD1	1:A:231:VAL:CG1	2.29	0.79
1:B:210:LEU:HD23	1:B:211:PHE:N	1.98	0.79
1:C:67:LEU:CD1	1:C:77:VAL:HB	2.11	0.79
1:B:210:LEU:HD11	1:B:405:VAL:CG1	2.13	0.78
1:C:217:GLY:CA	1:C:231:VAL:HG13	2.14	0.78
1:A:426:PHE:CD2	1:A:427:ARG:HG3	2.18	0.78
1:A:183:ALA:HB3	1:A:187:ILE:HD11	1.66	0.77
1:C:226:GLU:HB3	1:C:227:VAL:HA	1.67	0.76
1:C:372:ASP:HB3	1:C:374:ALA:O	1.85	0.76
1:C:66:ASN:HD21	1:C:235:MET:N	1.83	0.75
1:A:424:ASP:HB3	1:A:427:ARG:O	1.86	0.75
1:C:174:SER:HB2	1:C:176:TRP:CD1	2.20	0.75
1:C:67:LEU:HD11	1:C:77:VAL:CB	2.10	0.75
1:B:222:LEU:HD13	1:B:227:VAL:HA	1.67	0.74
1:A:97:SER:OG	1:A:187:ILE:CD1	2.35	0.74
1:C:123:ASP:OD2	1:C:125:ARG:HD3	1.87	0.74
1:A:235:MET:HE3	1:A:237:ILE:HD11	1.69	0.74
1:C:70:LYS:HG3	1:C:73:GLN:H	1.52	0.73
1:C:352:VAL:HB	1:C:355:GLN:HG2	1.71	0.73
1:C:70:LYS:HD2	1:C:73:GLN:HG3	1.72	0.72
1:A:339:ASN:OD1	1:A:340:ILE:HG23	1.89	0.71
1:B:274:LEU:O	1:B:275:LYS:HB2	1.90	0.71
1:C:351:GLU:H	1:C:355:GLN:NE2	1.89	0.71
1:C:229:ALA:CA	1:C:230:SER:CB	2.38	0.71
1:A:125:ARG:HG3	1:A:125:ARG:NH1	2.01	0.71

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:TRP:HE3	1:A:178:GLY:H	1.38	0.71
1:C:228:LEU:O	1:C:230:SER:HB3	1.90	0.70
1:A:97:SER:OG	1:A:187:ILE:CG1	2.39	0.70
1:C:337:PRO:HB2	1:C:340:ILE:CD1	2.22	0.70
1:B:294:ASN:ND2	1:B:384:ALA:O	2.25	0.70
1:A:97:SER:OG	1:A:187:ILE:HG13	1.93	0.69
1:B:310:LYS:CE	1:B:323:TRP:CD1	2.74	0.69
1:C:67:LEU:O	1:C:233:GLY:N	2.22	0.69
1:A:315:THR:HG23	1:B:339:ASN:HD21	1.58	0.68
1:B:100:ALA:HB2	1:B:161:ALA:HB3	1.75	0.68
1:A:254:ILE:O	1:A:411:LYS:HE2	1.95	0.67
1:C:337:PRO:HB2	1:C:340:ILE:HD12	1.76	0.67
1:B:210:LEU:HD23	1:B:210:LEU:C	2.14	0.67
1:C:69:GLY:HA3	1:C:75:TYR:HA	1.76	0.67
1:A:220:PHE:CE2	1:A:379:ASP:OD1	2.48	0.67
1:C:71:SER:HB2	1:C:400:GLU:OE1	1.96	0.66
1:A:68:ARG:O	1:A:76:TYR:CD1	2.49	0.66
1:B:294:ASN:CB	1:B:384:ALA:O	2.44	0.66
1:A:220:PHE:HE2	1:A:379:ASP:OD1	1.78	0.65
1:A:154:VAL:CG2	1:A:205:THR:HG21	2.26	0.65
1:A:292:THR:HA	2:D:2:ILE:O	1.95	0.65
1:A:70:LYS:HZ1	1:A:222:LEU:HD13	1.62	0.65
1:C:218:ALA:HB1	1:C:220:PHE:O	1.96	0.65
1:A:175:ASN:HB3	1:A:228:LEU:HD22	1.79	0.64
1:B:255:ARG:HE	1:B:440:MET:HE3	1.61	0.64
1:B:210:LEU:HD23	1:B:211:PHE:CA	2.26	0.64
1:C:210:LEU:CD1	1:C:239:GLY:N	2.60	0.64
1:C:318:PHE:HB3	1:C:319:PRO:HD2	1.79	0.64
1:C:66:ASN:OD1	1:C:67:LEU:N	2.30	0.64
1:A:148:ILE:HB	1:A:151:GLY:HA3	1.78	0.64
1:A:68:ARG:HD2	1:A:228:LEU:O	1.99	0.63
1:A:68:ARG:O	1:A:76:TYR:CE1	2.51	0.63
1:C:123:ASP:CG	1:C:125:ARG:HB2	2.19	0.63
1:A:154:VAL:HG21	1:A:205:THR:HG21	1.79	0.62
1:B:354:ASN:CB	1:B:436:VAL:HG12	2.29	0.62
1:A:295:LEU:HD22	1:A:398:ILE:HD11	1.81	0.62
1:A:222:LEU:O	1:A:222:LEU:HD12	2.00	0.62
1:C:319:PRO:O	1:C:322:PHE:HB3	2.00	0.62
1:C:70:LYS:HE3	1:C:224:GLN:HB2	1.81	0.61
1:A:210:LEU:HD11	1:A:239:GLY:CA	2.30	0.61
1:B:227:VAL:HG23	1:B:228:LEU:HD23	1.82	0.61

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:176:TRP:HE3	1:A:178:GLY:N	1.98	0.60
1:A:328:LEU:HD11	1:A:370:VAL:HG21	1.82	0.60
1:C:222:LEU:HD23	1:C:222:LEU:N	2.17	0.60
1:C:76:TYR:CE1	1:C:224:GLN:NE2	2.52	0.60
1:B:332:GLN:HA	1:B:378:ASP:OD1	2.01	0.60
1:C:289:ASP:OD2	2:F:4:36D:O2	2.20	0.60
1:A:436:VAL:HB	1:B:436:VAL:HG13	1.84	0.59
1:A:223:ASN:O	1:A:227:VAL:HG23	2.01	0.59
1:A:175:ASN:CB	1:A:228:LEU:HD22	2.32	0.59
1:B:83:SER:HB2	1:B:119:SER:HB3	1.84	0.59
1:B:68:ARG:NH1	1:B:228:LEU:HD22	2.17	0.59
1:C:305:ALA:O	1:C:309:ILE:CG1	2.50	0.59
1:B:319:PRO:O	1:B:322:PHE:HB3	2.03	0.58
1:C:66:ASN:OD1	1:C:67:LEU:HD13	2.02	0.58
1:B:100:ALA:CB	1:B:161:ALA:HB3	2.34	0.58
1:B:263:ILE:HG23	1:B:443:CYS:O	2.04	0.58
1:B:353:THR:OG1	1:B:437:THR:O	2.21	0.57
1:A:184:TYR:O	1:A:187:ILE:HG12	2.04	0.57
1:B:258:TRP:HD1	1:B:258:TRP:H	1.52	0.57
1:B:134:GLN:CD	2:E:3:TIH:HD	2.22	0.57
1:B:96:SER:O	1:B:183:ALA:HB3	2.05	0.57
1:C:132:TYR:CG	2:F:4:36D:H10	2.40	0.57
1:B:352:VAL:HG12	1:B:354:ASN:OD1	2.05	0.57
1:C:174:SER:HB2	1:C:176:TRP:NE1	2.19	0.57
1:A:360:THR:O	1:A:429:ALA:HB1	2.05	0.56
1:A:154:VAL:HG23	1:A:154:VAL:O	2.05	0.56
1:A:339:ASN:HD21	1:B:315:THR:HG23	1.70	0.56
1:C:210:LEU:CD1	1:C:239:GLY:CA	2.80	0.56
1:A:240:ILE:HG23	1:A:403:TYR:HE2	1.71	0.56
1:B:71:SER:HB3	1:B:400:GLU:OE1	2.06	0.55
1:C:67:LEU:CD1	1:C:77:VAL:CB	2.80	0.55
1:A:210:LEU:HD11	1:A:239:GLY:HA2	1.88	0.55
1:B:383:PHE:CZ	1:B:385:ILE:HB	2.41	0.55
1:A:175:ASN:HD22	1:A:228:LEU:HD22	1.72	0.55
1:B:210:LEU:HD12	1:B:407:ASP:HA	1.89	0.55
1:A:218:ALA:HB1	1:A:369:PRO:HG2	1.89	0.55
1:A:174:SER:HB2	1:A:176:TRP:CD1	2.41	0.55
1:B:294:ASN:CG	1:B:384:ALA:O	2.45	0.55
1:C:329:VAL:O	1:C:380:CYS:HA	2.07	0.55
1:B:310:LYS:HE2	1:B:323:TRP:NE1	2.23	0.54
1:A:66:ASN:HB2	1:A:150:HIS:CG	2.42	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:329:VAL:O	1:A:380:CYS:HA	2.07	0.54
1:C:123:ASP:OD1	1:C:125:ARG:HG3	2.06	0.54
1:A:205:THR:OG1	1:A:207:VAL:HG23	2.08	0.54
1:B:310:LYS:CE	1:B:323:TRP:NE1	2.71	0.54
1:A:70:LYS:NZ	1:A:222:LEU:CD1	2.56	0.54
1:A:199:ASP:OD1	1:A:408:ARG:NH2	2.41	0.53
1:B:277:ASP:OD1	1:B:278:CYS:N	2.41	0.53
1:A:75:TYR:OH	1:A:400:GLU:OE1	2.20	0.53
1:C:70:LYS:O	1:C:71:SER:C	2.46	0.53
1:C:289:ASP:O	1:C:395:GLY:HA2	2.09	0.53
1:B:289:ASP:OD2	2:E:4:36D:O2	2.27	0.53
1:C:124:LEU:N	1:C:125:ARG:HA	2.23	0.53
1:C:277:ASP:OD1	1:C:278:CYS:N	2.42	0.52
1:C:69:GLY:O	1:C:231:VAL:CB	2.27	0.52
1:C:293:THR:OG1	2:F:2:ILE:HG13	2.08	0.52
1:B:216:CYS:SG	1:B:420:CYS:CB	2.98	0.52
1:C:291:GLY:C	2:F:2:ILE:HD12	2.30	0.52
1:B:210:LEU:HD11	1:B:405:VAL:HG13	1.90	0.51
1:C:319:PRO:HB2	3:C:506:HOH:O	2.10	0.51
1:A:148:ILE:HG22	1:A:151:GLY:H	1.76	0.51
1:C:221:PRO:O	1:C:222:LEU:HD23	2.10	0.51
1:B:258:TRP:CG	1:B:259:TYR:N	2.77	0.51
1:C:78:GLU:O	1:C:149:PRO:HD2	2.11	0.51
1:C:266:ARG:HH11	1:C:266:ARG:HG2	1.76	0.51
1:A:258:TRP:HD1	1:A:258:TRP:H	1.59	0.51
1:A:339:ASN:OD1	1:A:340:ILE:N	2.43	0.51
1:A:93:ASP:OD1	2:D:4:36D:O1	2.29	0.51
1:C:169:PHE:CD1	2:F:4:36D:H9	2.46	0.51
1:C:364:GLN:OE1	1:C:364:GLN:N	2.43	0.51
1:B:441:GLU:OE1	1:C:273:ASP:OD2	2.28	0.51
1:C:174:SER:O	1:C:175:ASN:HB3	2.11	0.51
1:A:260:TYR:HB3	1:A:413:ILE:HD11	1.93	0.50
1:C:249:LEU:HD23	1:C:416:ALA:HB2	1.92	0.50
1:B:169:PHE:HD1	2:E:4:36D:H9	1.70	0.50
1:C:66:ASN:CG	1:C:67:LEU:HD13	2.31	0.50
1:C:296:ARG:HB3	1:C:388:SER:HB2	1.91	0.50
1:A:293:THR:OG1	2:D:2:ILE:HG12	2.12	0.50
1:A:201:LEU:HD11	1:A:207:VAL:HG21	1.94	0.50
1:C:226:GLU:HA	1:C:227:VAL:C	2.32	0.50
1:B:268:GLU:HG2	1:B:273:ASP:HA	1.94	0.50
1:A:135:GLY:HA2	1:A:167:ASP:O	2.12	0.49

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:103:ALA:CB	1:B:162:ALA:HB1	2.42	0.49
1:B:363:PRO:HA	1:B:366:TYR:CE2	2.47	0.49
1:B:90:ILE:HG21	1:B:180:LEU:HB2	1.94	0.49
1:C:124:LEU:HD12	1:C:141:LEU:HB3	1.95	0.49
1:C:124:LEU:HD12	1:C:141:LEU:CB	2.43	0.49
1:C:62:GLU:HG3	1:C:63:MET:HG3	1.95	0.49
1:C:66:ASN:OD1	1:C:67:LEU:HB2	2.13	0.49
1:C:210:LEU:HD11	1:C:238:GLY:C	2.34	0.49
1:C:68:ARG:O	1:C:76:TYR:CE1	2.66	0.49
1:C:216:CYS:O	1:C:231:VAL:CG1	2.56	0.48
1:B:377:GLN:HA	1:B:377:GLN:NE2	2.28	0.48
1:A:158:ALA:O	1:A:160:ILE:HD12	2.13	0.48
1:A:235:MET:CE	1:A:237:ILE:CD1	2.77	0.48
1:B:176:TRP:CE3	1:B:178:GLY:O	2.66	0.48
1:C:333:ALA:HB2	1:C:377:GLN:O	2.12	0.48
1:A:426:PHE:CE2	1:A:427:ARG:HG3	2.49	0.48
1:C:352:VAL:CB	1:C:355:GLN:HG2	2.43	0.48
1:C:294:ASN:ND2	1:C:386:SER:OG	2.41	0.48
1:B:185:ALA:O	1:B:186:GLU:C	2.53	0.47
1:B:72:GLY:O	2:E:1:GLU:N	2.40	0.47
1:A:322:PHE:CZ	1:A:383:PHE:HB2	2.49	0.47
1:B:331:TRP:CE3	1:B:336:THR:HG23	2.50	0.47
1:C:226:GLU:CB	1:C:227:VAL:HA	2.37	0.47
1:A:363:PRO:HA	1:A:366:TYR:CE2	2.50	0.47
1:C:124:LEU:HB2	1:C:125:ARG:HA	1.96	0.47
1:A:151:GLY:O	1:A:152:PRO:C	2.54	0.47
1:B:274:LEU:O	1:B:275:LYS:CB	2.61	0.47
1:B:352:VAL:CG1	1:B:354:ASN:OD1	2.62	0.47
1:B:118:SER:HB3	1:B:121:TYR:HB2	1.97	0.47
1:A:222:LEU:C	1:A:222:LEU:HD12	2.35	0.46
1:A:183:ALA:CB	1:A:187:ILE:HD11	2.41	0.46
1:A:293:THR:O	1:A:397:VAL:HG13	2.15	0.46
1:B:440:MET:O	1:B:442:ASP:O	2.33	0.46
1:A:310:LYS:HG2	1:A:323:TRP:CE2	2.51	0.46
1:A:426:PHE:CG	1:B:311:ALA:HB1	2.51	0.46
1:C:369:PRO:O	1:C:370:VAL:HG13	2.15	0.46
1:B:299:LYS:HB2	1:B:387:GLN:HE21	1.80	0.46
1:B:192:ASP:OD1	1:B:193:SER:N	2.49	0.46
1:B:134:GLN:HB2	2:E:3:TIH:HB3	1.96	0.46
1:B:169:PHE:CE1	2:E:4:36D:C5	2.99	0.46
1:A:169:PHE:CD1	2:D:4:36D:H9	2.51	0.46

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:220:PHE:CD1	1:A:221:PRO:HA	2.51	0.46
1:B:176:TRP:HE3	1:B:178:GLY:O	1.98	0.46
1:A:250:TRP:CZ3	1:C:352:VAL:HG13	2.50	0.46
1:B:376:SER:O	1:B:377:GLN:C	2.53	0.46
1:B:169:PHE:CE1	2:E:4:36D:H9	2.50	0.46
1:A:405:VAL:O	1:A:413:ILE:HA	2.16	0.46
1:A:60:PHE:CZ	1:A:239:GLY:HA3	2.51	0.45
1:B:354:ASN:HA	1:B:436:VAL:HA	1.98	0.45
1:B:175:ASN:O	1:B:175:ASN:CG	2.55	0.45
1:B:354:ASN:HB3	1:B:436:VAL:HG12	1.98	0.45
1:A:245:TYR:CD1	1:A:403:TYR:CD2	3.04	0.45
1:C:291:GLY:O	2:F:2:ILE:HD12	2.16	0.45
1:A:426:PHE:CE1	1:B:311:ALA:HB1	2.50	0.45
1:C:401:GLY:O	1:C:418:SER:HB3	2.16	0.45
1:B:310:LYS:HE3	1:B:323:TRP:CG	2.48	0.45
1:B:134:GLN:OE1	2:E:2:ILE:HA	2.15	0.45
1:B:310:LYS:HG3	1:B:323:TRP:CZ2	2.51	0.45
1:B:347:TYR:CZ	1:B:358:ARG:HD2	2.52	0.45
1:C:306:VAL:HA	1:C:309:ILE:HD12	1.99	0.45
1:C:68:ARG:HA	1:C:232:GLY:HA2	1.98	0.45
1:A:132:TYR:CG	2:D:4:36D:H10	2.52	0.45
1:B:132:TYR:CG	2:E:4:36D:H10	2.52	0.45
1:A:176:TRP:CE3	1:A:178:GLY:N	2.84	0.45
1:A:313:SER:HB2	1:A:340:ILE:HD12	1.98	0.45
1:C:305:ALA:O	1:C:309:ILE:CD1	2.65	0.45
1:C:230:SER:C	1:C:231:VAL:HG23	2.37	0.45
1:B:210:LEU:CD2	1:B:211:PHE:N	2.75	0.44
1:A:76:TYR:OH	1:A:228:LEU:CD2	2.65	0.44
1:C:265:VAL:HG23	1:C:347:TYR:HB2	1.98	0.44
1:A:210:LEU:C	1:A:210:LEU:HD12	2.37	0.44
1:A:227:VAL:O	1:A:230:SER:O	2.36	0.44
1:A:82:GLY:HA2	1:A:144:ASP:OD1	2.18	0.44
1:B:60:PHE:CZ	1:B:239:GLY:HA3	2.52	0.44
1:C:362:LEU:HB3	1:C:363:PRO:HD2	2.00	0.44
1:B:106:HIS:CG	1:B:107:PRO:HD2	2.53	0.44
1:C:384:ALA:CB	1:C:397:VAL:HG11	2.48	0.43
1:B:169:PHE:CD1	2:E:4:36D:C5	2.94	0.43
1:C:123:ASP:OD1	1:C:125:ARG:HB2	2.18	0.43
1:A:89:ASN:HB2	1:A:176:TRP:HA	2.00	0.43
1:C:116:GLN:HG2	1:C:117:LEU:HD12	1.98	0.43
1:C:124:LEU:CD1	1:C:141:LEU:HB3	2.48	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:174:SER:O	1:C:175:ASN:CB	2.65	0.43
1:C:134:GLN:HB2	2:F:3:TIH:HB3	1.99	0.43
1:C:267:VAL:CG2	1:C:281:TYR:CE1	3.02	0.43
1:A:215:LEU:O	1:A:400:GLU:HA	2.18	0.43
1:A:220:PHE:HA	1:A:221:PRO:HA	1.70	0.43
1:B:79:MET:SD	1:B:90:ILE:HG13	2.58	0.43
1:C:113:TYR:OH	1:C:144:ASP:OD2	2.21	0.43
1:C:229:ALA:CB	1:C:230:SER:OG	2.60	0.43
1:C:68:ARG:HH11	1:C:232:GLY:HA3	1.84	0.43
1:C:294:ASN:ND2	2:F:1:GLU:OE2	2.52	0.43
1:A:65:ASP:CB	1:A:231:VAL:HG11	2.49	0.43
1:A:313:SER:CB	1:A:340:ILE:HD12	2.49	0.43
1:B:302:PHE:CE1	1:B:385:ILE:HG22	2.54	0.42
1:C:265:VAL:HG23	1:C:266:ARG:N	2.34	0.42
1:B:289:ASP:OD1	1:B:291:GLY:N	2.24	0.42
1:A:341:PHE:HB3	1:A:363:PRO:HB3	2.01	0.42
1:C:306:VAL:HG23	1:C:323:TRP:CE3	2.54	0.42
1:A:372:ASP:OD1	1:A:375:THR:HA	2.19	0.42
1:B:174:SER:O	1:B:175:ASN:HB3	2.20	0.42
1:B:259:TYR:O	1:B:261:GLU:N	2.53	0.42
1:B:354:ASN:HB2	1:B:436:VAL:HG12	2.02	0.42
1:C:345:SER:HA	1:C:359:ILE:O	2.20	0.42
1:C:369:PRO:O	1:C:370:VAL:CG1	2.68	0.42
1:B:214:GLN:NE2	1:B:420:CYS:SG	2.93	0.42
1:C:377:GLN:HA	1:C:377:GLN:NE2	2.34	0.42
1:C:283:TYR:HA	1:C:284:ASP:HA	1.79	0.42
1:C:319:PRO:HB2	1:C:327:GLN:OE1	2.20	0.42
1:C:78:GLU:HG2	1:C:149:PRO:HG3	2.01	0.42
1:B:192:ASP:O	1:B:193:SER:C	2.58	0.42
1:B:103:ALA:HB2	1:B:162:ALA:HB1	2.02	0.41
1:C:222:LEU:O	1:C:230:SER:N	2.45	0.41
1:C:363:PRO:HA	1:C:366:TYR:CE1	2.55	0.41
1:C:338:TRP:HZ3	1:C:367:LEU:HD12	1.85	0.41
1:B:101:VAL:O	1:B:101:VAL:HG23	2.20	0.41
1:B:96:SER:C	1:B:183:ALA:HB3	2.40	0.41
1:C:260:TYR:HB3	1:C:413:ILE:HD11	2.01	0.41
1:B:199:ASP:OD1	1:B:408:ARG:NH2	2.52	0.41
1:B:374:ALA:O	1:B:375:THR:C	2.59	0.41
1:C:256:ARG:HD3	1:C:258:TRP:NE1	2.35	0.41
1:A:241:ASP:OD1	1:A:243:SER:OG	2.33	0.41
1:B:175:ASN:O	1:B:175:ASN:OD1	2.39	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:70:LYS:C	1:C:72:GLY:N	2.73	0.41
1:A:66:ASN:CB	1:A:150:HIS:CG	3.04	0.41
1:B:80:THR:O	1:B:146:VAL:HA	2.20	0.41
1:A:259:TYR:CE1	2:D:4:36D:H19	2.55	0.41
1:A:258:TRP:N	1:A:258:TRP:CD1	2.89	0.41
1:C:267:VAL:HG23	1:C:281:TYR:CE1	2.55	0.41
1:A:97:SER:HG	1:A:187:ILE:CG1	2.33	0.41
1:A:225:SER:HA	1:A:228:LEU:HD12	2.03	0.41
1:C:297:LEU:O	1:C:388:SER:N	2.54	0.41
1:C:376:SER:HB3	1:C:378:ASP:CG	2.40	0.41
1:A:196:PRO:O	1:A:197:PHE:C	2.60	0.40
1:A:76:TYR:OH	1:A:228:LEU:HD22	2.21	0.40
1:A:318:PHE:HD1	1:A:329:VAL:HG11	1.86	0.40
1:B:143:THR:OG1	1:B:157:ARG:NH1	2.54	0.40
1:B:134:GLN:OE1	2:E:3:TIH:CD	2.35	0.40
1:A:289:ASP:OD2	2:D:4:36D:O2	2.40	0.40
1:B:283:TYR:HA	1:B:284:ASP:HA	1.85	0.40
1:C:68:ARG:O	1:C:76:TYR:CD1	2.75	0.40
1:C:171:ILE:HB	1:C:174:SER:HB3	2.02	0.40
1:C:269:ILE:CG1	1:C:274:LEU:HD21	2.52	0.40
1:C:63:MET:HE1	1:C:152:PRO:HD3	2.03	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	385/388 (99%)	381 (99%)	3 (1%)	1 (0%)	44	75
1	B	386/388 (100%)	378 (98%)	7 (2%)	1 (0%)	44	75
1	C	386/388 (100%)	379 (98%)	6 (2%)	1 (0%)	44	75
2	D	1/4 (25%)	1 (100%)	0	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	E	1/4 (25%)	1 (100%)	0	0	100	100
2	F	1/4 (25%)	1 (100%)	0	0	100	100
All	All	1160/1176 (99%)	1141 (98%)	16 (1%)	3 (0%)	44	75

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	131	PRO
1	B	131	PRO
1	C	131	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	330/331 (100%)	328 (99%)	2 (1%)	89	96
1	B	331/331 (100%)	329 (99%)	2 (1%)	89	96
1	C	331/331 (100%)	325 (98%)	6 (2%)	64	87
2	D	2/2 (100%)	2 (100%)	0	100	100
2	E	2/2 (100%)	2 (100%)	0	100	100
2	F	2/2 (100%)	1 (50%)	1 (50%)	0	0
All	All	998/999 (100%)	987 (99%)	11 (1%)	78	92

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	216	CYS
1	A	258	TRP
1	B	228	LEU
1	B	258	TRP
1	C	216	CYS
1	C	258	TRP
1	C	266	ARG

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	C	278	CYS
1	C	352	VAL
1	C	355	GLN
2	F	2	ILE

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

Mol	Chain	Res	Type
1	A	175	ASN
1	A	365	GLN
1	B	175	ASN
1	B	339	ASN
1	B	365	GLN
1	B	377	GLN
1	B	387	GLN
1	C	134	GLN
1	C	223	ASN
1	C	332	GLN
1	C	355	GLN
1	C	377	GLN
1	C	387	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

3 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	$\# Z > 2$
2	TIH	D	3	2	8,10,11	1.33	1 (12%)	8,12,14	2.37	2 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	TIH	E	3	2	8,10,11	1.06	1 (12%)	8,12,14	2.03	2 (25%)
2	TIH	F	3	2	8,10,11	1.21	1 (12%)	8,12,14	2.51	3 (37%)

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
2	TIH	D	3	2	-	0/3/6/8	0/1/1/1
2	TIH	E	3	2	-	0/3/6/8	0/1/1/1
2	TIH	F	3	2	-	0/3/6/8	0/1/1/1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	3	TIH	CE1-CD	-2.23	1.32	1.39
2	F	3	TIH	CA-C	2.36	1.53	1.50
2	D	3	TIH	CA-C	2.90	1.54	1.50

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	3	TIH	CE1-CE2-SD	-5.32	107.45	113.18
2	F	3	TIH	CE1-CE2-SD	-4.79	108.03	113.18
2	E	3	TIH	CE1-CE2-SD	-4.28	108.57	113.18
2	F	3	TIH	CB-CA-C	-3.45	104.76	111.41
2	D	3	TIH	CB-CA-C	-3.04	105.55	111.41
2	E	3	TIH	CB-CA-C	-2.99	105.66	111.41
2	F	3	TIH	CE2-SD-CG	3.22	93.72	91.95

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	3	TIH	4	0
2	F	3	TIH	1	0

5.5 Carbohydrates [i](#)

There are no carbohydrates 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

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	387/388 (99%)	0.38	26 (6%) 19 14	35, 55, 120, 207	0
1	B	388/388 (100%)	0.28	26 (6%) 19 14	33, 56, 106, 189	0
1	C	388/388 (100%)	0.47	38 (9%) 8 5	34, 61, 134, 223	0
2	D	2/4 (50%)	0.51	0 100 100	49, 49, 49, 61	0
2	E	2/4 (50%)	0.12	0 100 100	47, 47, 47, 54	0
2	F	2/4 (50%)	0.42	0 100 100	60, 60, 60, 65	0
All	All	1169/1176 (99%)	0.37	90 (7%) 14 10	33, 57, 123, 223	0

All (90) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	221	PRO	15.7
1	C	223	ASN	14.9
1	A	220	PHE	14.9
1	B	229	ALA	12.1
1	A	222	LEU	11.6
1	A	219	GLY	11.1
1	C	221	PRO	10.8
1	B	220	PHE	10.5
1	C	224	GLN	10.1
1	A	226	GLU	10.0
1	B	221	PRO	9.5
1	C	229	ALA	9.5
1	A	230	SER	9.1
1	C	220	PHE	9.0
1	C	225	SER	8.8
1	A	229	ALA	8.8
1	B	228	LEU	8.8
1	C	219	GLY	8.5
1	C	373	VAL	7.7

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	222	LEU	7.7
1	A	373	VAL	7.3
1	C	375	THR	7.0
1	C	227	VAL	6.7
1	C	218	ALA	6.5
1	A	374	ALA	6.4
1	A	231	VAL	6.2
1	C	226	GLU	6.1
1	A	223	ASN	6.0
1	B	222	LEU	6.0
1	B	374	ALA	5.3
1	C	228	LEU	5.1
1	B	223	ASN	5.0
1	B	375	THR	5.0
1	B	224	GLN	4.8
1	B	225	SER	4.7
1	C	374	ALA	4.7
1	C	371	GLU	4.5
1	C	315	THR	4.4
1	A	228	LEU	4.2
1	A	154	VAL	4.1
1	C	124	LEU	3.9
1	A	218	ALA	3.8
1	C	230	SER	3.7
1	A	150	HIS	3.7
1	C	125	ARG	3.6
1	B	219	GLY	3.6
1	A	224	GLN	3.6
1	B	373	VAL	3.4
1	B	227	VAL	3.4
1	C	84	PRO	3.4
1	C	70	LYS	3.4
1	A	227	VAL	3.4
1	B	377	GLN	3.3
1	B	226	GLU	3.2
1	B	206	HIS	3.2
1	A	317	LYS	3.2
1	A	377	GLN	3.1
1	A	225	SER	3.1
1	A	425	GLU	3.1
1	C	377	GLN	2.9
1	A	375	THR	2.9

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	B	145	LEU	2.9
1	C	217	GLY	2.7
1	C	319	PRO	2.7
1	C	323	TRP	2.7
1	C	117	LEU	2.6
1	A	372	ASP	2.6
1	C	425	GLU	2.5
1	C	318	PHE	2.5
1	C	328	LEU	2.5
1	B	333	ALA	2.4
1	C	313	SER	2.3
1	B	329	VAL	2.3
1	C	372	ASP	2.3
1	A	151	GLY	2.3
1	B	218	ALA	2.2
1	C	322	PHE	2.2
1	C	426	PHE	2.2
1	C	233	GLY	2.2
1	B	317	LYS	2.1
1	B	105	PRO	2.1
1	C	68	ARG	2.1
1	B	230	SER	2.1
1	A	217	GLY	2.1
1	B	443	CYS	2.1
1	B	384	ALA	2.1
1	B	318	PHE	2.1
1	C	376	SER	2.1
1	A	430	ALA	2.0
1	C	380	CYS	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(Å ²)	Q<0.9
2	TIH	E	3	10/11	0.93	0.16	-	36,42,45,56	0
2	TIH	D	3	10/11	0.96	0.16	-	35,41,44,58	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
2	TIH	F	3	10/11	0.97	0.13	-	49,56,59,61	0

6.3 Carbohydrates [i](#)

There are no carbohydrates 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.