



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 02:40 AM UTC

PDB ID : 4TRW / pdb_00004trw
Title : Structure of BACE1 complex with a syn-HEA-type inhibitor
Authors : Akaji, K.; Teruya, K.; Akiyama, T.; Sanjho, A.; Yamashita, E.; Nakagawa, A.
Deposited on : 2014-06-18
Resolution : 2.85 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

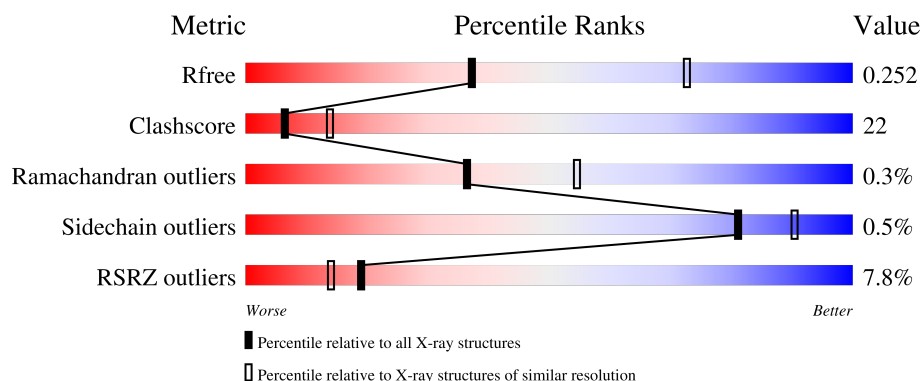
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.85 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	1407 (2.88-2.84)
Clashscore	190562	1446 (2.88-2.84)
Ramachandran outliers	187476	1406 (2.88-2.84)
Sidechain outliers	187428	1407 (2.88-2.84)
RSRZ outliers	180081	1408 (2.88-2.84)

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	390	6% 73% 25% .
1	B	390	8% 69% 27% .
1	C	390	9% 71% 26% .
2	D	4	50% 25% 25%
2	E	4	25% 50% 25%

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Mol	Chain	Length	Quality of chain
2	F	4	 A horizontal bar chart showing the quality of chain F. The bar is divided into three segments: a yellow segment on the left labeled '50%', an orange segment in the middle labeled '25%', and a red segment on the right labeled '25%'. The segments are separated by thin black lines.

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 9353 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	389	Total	C	N	O	S	0	0	0
			3056	1957	507	578	14			
1	B	389	Total	C	N	O	S	0	0	0
			3056	1957	507	578	14			
1	C	390	Total	C	N	O	S	0	0	0
			3060	1959	508	579	14			

- Molecule 2 is a protein called L-alpha-glutamyl-L-isoleucyl-N-[(2R,3S)-1-{[(1S)-1-carboxybutyl]amino}-2-hydroxy-5-methylhexan-3-yl]-3-thiophen-2-yl-L-alaninamide.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	4	Total	C	N	O	S	0	0	0
			44	30	5	8	1			
2	E	4	Total	C	N	O	S	0	0	0
			44	30	5	8	1			
2	F	4	Total	C	N	O	S	0	0	0
			44	30	5	8	1			

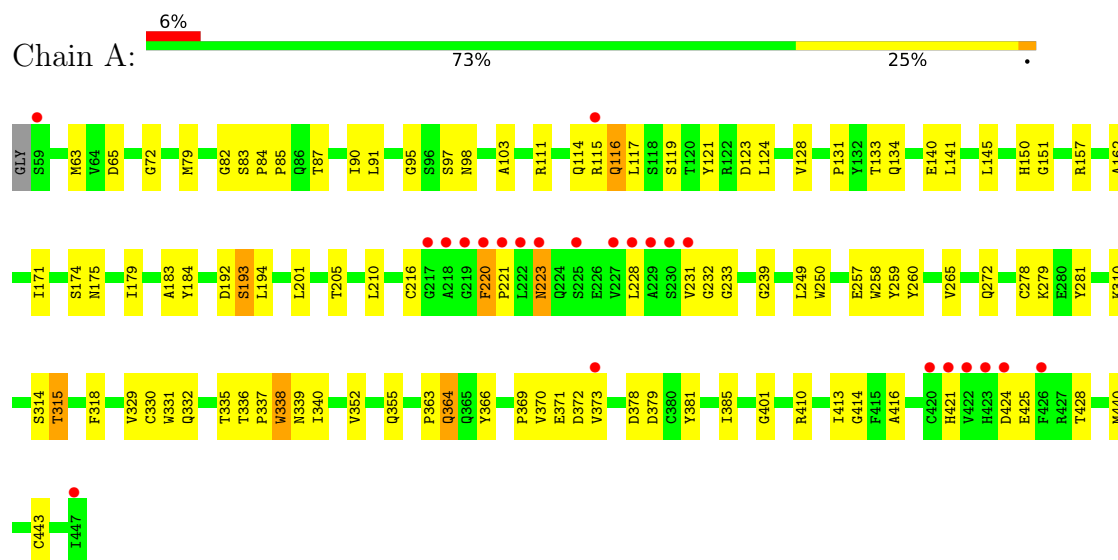
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	12	Total	O	0	0
			12	12		
3	B	21	Total	O	0	0
			21	21		
3	C	16	Total	O	0	0
			16	16		

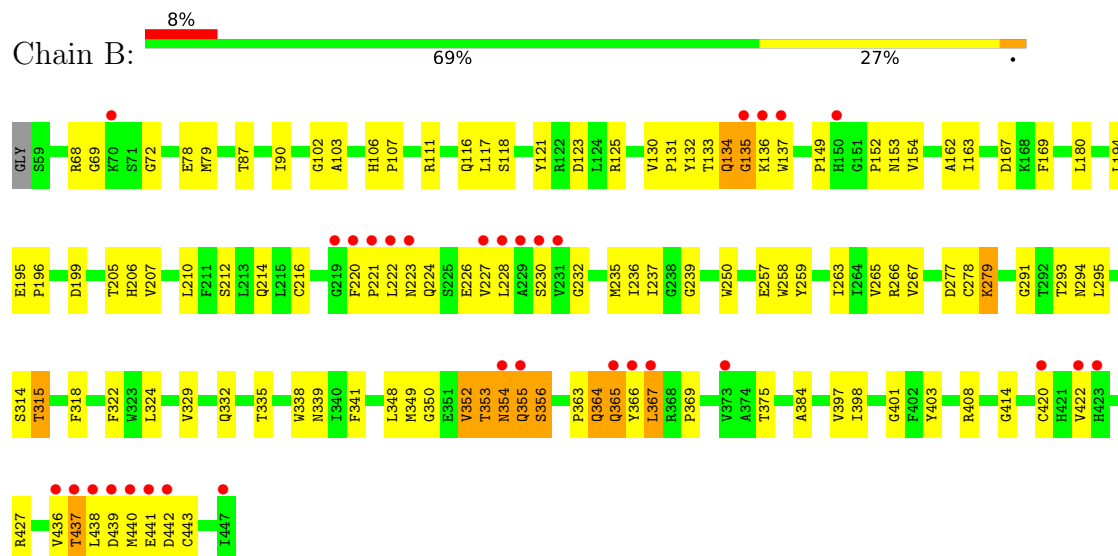
3 Residue-property plots

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: Beta-secretase 1

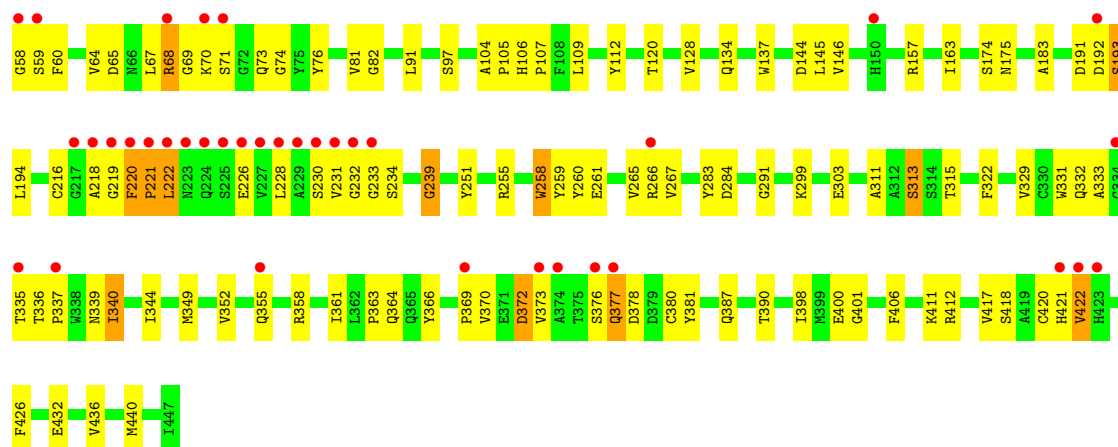


• Molecule 1: Beta-secretase 1



• Molecule 1: Beta-secretase 1





- Molecule 2: L-alpha-glutamyl-L-isoleucyl-N-[(2R,3S)-1-([(1S)-1-carboxybutyl]amino)-2-hydroxy-5-methylhexan-3-yl]-3-thiophen-2-yl-L-alaninamide

Chain D: 50% 25% 25%



- Molecule 2: L-alpha-glutamyl-L-isoleucyl-N-[(2R,3S)-1-([(1S)-1-carboxybutyl]amino)-2-hydroxy-5-methylhexan-3-yl]-3-thiophen-2-yl-L-alaninamide

Chain E: 25% 50% 25%



- Molecule 2: L-alpha-glutamyl-L-isoleucyl-N-[(2R,3S)-1-([(1S)-1-carboxybutyl]amino)-2-hydroxy-5-methylhexan-3-yl]-3-thiophen-2-yl-L-alaninamide

Chain F: 50% 25% 25%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	82.28Å 103.66Å 102.00Å 90.00° 102.73° 90.00°	Depositor
Resolution (Å)	49.72 – 2.85 49.72 – 2.85	Depositor EDS
% Data completeness (in resolution range)	99.1 (49.72-2.85) 99.1 (49.72-2.85)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.29 (at 2.86Å)	Xtriage
Refinement program	REFMAC 5.7.0032	Depositor
R, R_{free}	0.208 , 0.238 0.207 , 0.252	Depositor DCC
R_{free} test set	1950 reflections (4.98%)	wwPDB-VP
Wilson B-factor (Å ²)	61.4	Xtriage
Anisotropy	0.044	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 36.8	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	0.93	EDS
Total number of atoms	9353	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.42% 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: LHE, 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	0.80	0/3134	1.06	14/4262 (0.3%)
1	B	0.83	0/3134	1.12	21/4262 (0.5%)
1	C	0.79	0/3138	1.13	19/4267 (0.4%)
2	D	0.77	0/16	0.46	0/20
2	E	0.02	0/16	0.28	0/20
2	F	0.02	0/16	0.28	0/20
All	All	0.81	0/9454	1.10	54/12851 (0.4%)

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
2	D	0	1
2	E	0	1
2	F	0	1
All	All	0	3

There are no bond length outliers.

All (54) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	220	PHE	C-N-CD	-13.29	91.37	120.60
1	C	426	PHE	N-CA-C	11.30	123.35	111.14
1	C	234	SER	N-CA-C	10.16	125.08	108.52
1	C	69	GLY	N-CA-C	9.65	122.73	111.63
1	B	365	GLN	N-CA-C	9.37	121.25	111.14
1	C	311	ALA	N-CA-C	9.10	120.97	111.14
1	A	233	GLY	N-CA-C	8.64	122.58	110.74

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	364	GLN	N-CA-C	-8.28	102.25	111.28
1	A	257	GLU	N-CA-C	8.13	123.31	110.70
1	B	224	GLN	N-CA-C	7.42	121.25	111.75
1	A	116	GLN	N-CA-C	7.20	121.76	113.19
1	C	68	ARG	N-CA-C	7.18	121.36	109.95
1	B	257	GLU	N-CA-C	7.03	120.45	110.68
1	C	193	SER	N-CA-C	-6.85	103.83	111.71
1	B	135	GLY	N-CA-C	6.77	123.03	111.04
1	B	134	GLN	CA-C-N	-6.40	117.66	122.33
1	B	134	GLN	C-N-CA	-6.40	117.66	122.33
1	B	355	GLN	N-CA-C	6.30	118.14	111.28
1	C	239	GLY	N-CA-C	6.27	119.77	110.42
1	B	279	LYS	N-CA-C	-6.09	104.64	111.28
1	C	112	TYR	N-CA-C	6.05	117.44	108.60
1	B	367	LEU	CA-C-N	-6.04	114.46	122.56
1	B	367	LEU	C-N-CA	-6.04	114.46	122.56
1	B	133	THR	N-CA-C	-6.02	104.72	111.28
1	A	220	PHE	N-CA-C	-5.98	99.17	109.15
1	A	98	ASN	N-CA-C	5.92	119.28	109.46
1	B	353	THR	N-CA-C	-5.89	105.19	112.38
1	C	267	VAL	N-CA-CB	-5.78	101.24	111.93
1	C	313	SER	CB-CA-C	-5.76	100.46	110.09
1	B	354	ASN	N-CA-C	-5.68	105.17	111.36
1	A	310	LYS	N-CA-C	-5.65	104.81	110.97
1	B	352	VAL	N-CA-C	-5.63	99.66	108.95
1	C	68	ARG	N-CA-CB	-5.63	101.84	110.85
1	C	265	VAL	N-CA-C	5.59	119.52	113.43
1	C	222	LEU	N-CA-C	5.52	118.45	110.28
1	A	128	VAL	N-CA-C	5.50	115.60	108.35
1	A	338	TRP	N-CA-C	5.49	117.97	111.33
1	A	315	THR	N-CA-C	-5.47	105.32	111.28
1	C	372	ASP	N-CA-C	5.46	118.54	110.46
1	C	303	GLU	N-CA-C	5.42	117.27	111.36
1	B	257	GLU	CB-CA-C	-5.42	104.24	111.88
1	B	230	SER	N-CA-C	-5.41	102.04	110.10
1	A	98	ASN	N-CA-CB	-5.41	101.23	110.16
1	B	315	THR	N-CA-C	-5.41	105.39	111.28
1	C	398	ILE	CB-CA-C	-5.40	104.95	112.02
1	C	390	THR	N-CA-C	5.27	119.26	112.41
1	A	364	GLN	N-CA-C	-5.25	105.56	111.28
1	A	193	SER	N-CA-C	-5.23	105.54	112.34
1	A	223	ASN	N-CA-C	5.07	117.39	110.55

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	220	PHE	CA-C-N	-5.06	114.57	119.78
1	B	220	PHE	C-N-CA	-5.06	114.57	119.78
1	B	356	SER	N-CA-C	5.03	117.33	110.24
1	A	265	VAL	CB-CA-C	-5.02	104.97	111.20
1	C	422	VAL	N-CA-C	-5.02	102.41	109.63

There are no chirality outliers.

All (3) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	3	TIH	Peptide
2	E	3	TIH	Peptide
2	F	3	TIH	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	3056	0	2969	100	0
1	B	3056	0	2969	168	0
1	C	3060	0	2970	145	0
2	D	44	0	39	5	0
2	E	44	0	40	6	0
2	F	44	0	40	5	0
3	A	12	0	0	2	0
3	B	21	0	0	1	0
3	C	16	0	0	1	0
All	All	9353	0	9027	399	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 (399) 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:219:GLY:HA3	1:C:220:PHE:CD2	1.42	1.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:363:PRO:HA	1:B:366:TYR:CE2	1.47	1.49
1:B:263:ILE:HG21	1:B:440:MET:CE	1.59	1.31
1:A:352:VAL:HG11	1:C:355:GLN:NE2	1.54	1.22
1:A:338:TRP:CZ3	1:A:364:GLN:HA	1.77	1.18
1:C:369:PRO:HG3	1:C:381:TYR:CE1	1.79	1.16
1:B:263:ILE:CG2	1:B:440:MET:CE	2.25	1.15
1:B:436:VAL:HG21	1:C:436:VAL:CG2	1.78	1.14
1:C:219:GLY:HA3	1:C:220:PHE:CE2	1.82	1.14
1:B:263:ILE:CG2	1:B:440:MET:HE1	1.78	1.13
1:B:265:VAL:HG21	1:B:437:THR:OG1	1.46	1.12
1:B:356:SER:HB3	1:B:437:THR:HG21	1.27	1.11
1:B:436:VAL:CG2	1:C:436:VAL:HG23	1.79	1.11
1:C:219:GLY:CA	1:C:220:PHE:CD2	2.33	1.10
1:C:331:TRP:HE3	1:C:336:THR:CG2	1.65	1.08
1:C:352:VAL:HB	1:C:355:GLN:OE1	1.53	1.08
1:C:219:GLY:HA3	1:C:220:PHE:CG	1.88	1.08
1:B:363:PRO:HB3	1:B:366:TYR:OH	1.54	1.08
1:C:331:TRP:HB2	1:C:336:THR:CG2	1.83	1.08
1:C:331:TRP:HB2	1:C:336:THR:HG22	1.11	1.08
1:B:363:PRO:CA	1:B:366:TYR:CE2	2.36	1.07
1:A:352:VAL:HG11	1:C:355:GLN:HE22	0.93	1.07
1:A:338:TRP:CH2	1:A:364:GLN:HA	1.91	1.05
1:B:356:SER:HB3	1:B:437:THR:CG2	1.85	1.05
1:B:132:TYR:HE2	1:B:137:TRP:NE1	1.54	1.04
1:C:219:GLY:CA	1:C:220:PHE:CG	2.40	1.04
1:C:222:LEU:HD22	1:C:226:GLU:CD	1.84	1.03
1:C:331:TRP:CE3	1:C:336:THR:CG2	2.40	1.03
1:B:356:SER:CB	1:B:437:THR:CG2	2.36	1.03
1:B:440:MET:HB3	1:B:441:GLU:HA	1.36	1.02
1:C:331:TRP:CB	1:C:336:THR:HG22	1.90	0.99
1:C:231:VAL:H	1:C:232:GLY:HA2	1.27	0.99
1:B:366:TYR:O	1:B:384:ALA:HB3	1.62	0.98
1:B:263:ILE:HG21	1:B:440:MET:HE2	1.41	0.98
1:B:356:SER:OG	1:B:437:THR:HG22	1.65	0.97
1:C:220:PHE:HB3	1:C:369:PRO:HB2	1.46	0.97
1:B:278:CYS:CB	1:B:443:CYS:SG	2.53	0.97
1:C:220:PHE:CB	1:C:369:PRO:HB2	1.95	0.96
1:C:331:TRP:CE3	1:C:336:THR:HB	2.01	0.96
1:B:436:VAL:HG21	1:C:436:VAL:HG23	0.96	0.95
1:B:356:SER:CB	1:B:437:THR:HG21	1.96	0.95
1:A:352:VAL:CG1	1:C:355:GLN:HE22	1.80	0.94

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:223:ASN:OD1	1:B:226:GLU:HB2	1.66	0.94
1:C:331:TRP:CE3	1:C:336:THR:HG22	2.03	0.94
1:C:369:PRO:HG3	1:C:381:TYR:HE1	1.29	0.93
1:A:338:TRP:CZ3	1:A:364:GLN:CA	2.52	0.92
1:C:372:ASP:OD1	1:C:373:VAL:N	2.03	0.92
1:C:222:LEU:HD22	1:C:226:GLU:OE2	1.70	0.91
1:C:81:VAL:HG12	1:C:146:VAL:HG22	1.53	0.91
1:B:265:VAL:HG21	1:B:437:THR:CB	2.02	0.90
1:B:353:THR:HG23	1:B:437:THR:O	1.71	0.90
1:B:291:GLY:O	2:E:2:ILE:HD12	1.71	0.89
1:C:417:VAL:HG13	1:C:421:HIS:NE2	1.88	0.89
1:A:364:GLN:NE2	1:A:424:ASP:HB2	1.89	0.88
1:B:349:MET:SD	1:B:440:MET:SD	2.72	0.88
1:C:369:PRO:HG3	1:C:381:TYR:CD1	2.09	0.88
1:B:278:CYS:HB2	1:B:443:CYS:SG	2.14	0.88
1:B:279:LYS:HG2	1:B:442:ASP:O	1.72	0.87
1:B:439:ASP:OD2	1:C:266:ARG:HD2	1.74	0.87
1:B:363:PRO:HA	1:B:366:TYR:CZ	2.08	0.87
1:B:265:VAL:CG1	1:B:440:MET:HE3	2.04	0.87
1:B:265:VAL:CG2	1:B:437:THR:OG1	2.22	0.86
1:C:331:TRP:CE3	1:C:336:THR:CB	2.58	0.86
1:B:436:VAL:CG2	1:C:436:VAL:CG2	2.46	0.86
1:C:231:VAL:N	1:C:232:GLY:HA2	1.83	0.85
1:C:67:LEU:HB2	1:C:233:GLY:O	1.78	0.83
1:B:294:ASN:HB3	1:B:384:ALA:O	1.77	0.83
1:B:265:VAL:HG12	1:B:440:MET:HE3	1.60	0.83
1:B:440:MET:CB	1:B:441:GLU:HA	2.05	0.82
1:B:356:SER:CB	1:B:437:THR:HG22	2.08	0.82
1:B:366:TYR:CE1	1:B:367:LEU:HG	2.15	0.81
1:B:222:LEU:HD12	1:B:222:LEU:O	1.80	0.81
1:B:356:SER:OG	1:B:437:THR:CG2	2.26	0.81
1:B:263:ILE:CG2	1:B:440:MET:HE2	2.01	0.80
1:C:331:TRP:HE3	1:C:336:THR:HG21	1.47	0.80
1:C:355:GLN:O	1:C:440:MET:SD	2.40	0.80
1:C:231:VAL:HB	1:C:233:GLY:N	1.97	0.80
1:B:265:VAL:HG11	1:B:437:THR:OG1	1.81	0.80
1:C:220:PHE:CD2	1:C:369:PRO:HG2	2.17	0.79
1:B:440:MET:HA	1:B:442:ASP:H	1.47	0.78
1:A:318:PHE:HD1	1:A:329:VAL:HG11	1.47	0.78
1:B:366:TYR:O	1:B:384:ALA:CB	2.32	0.78
1:C:230:SER:OG	1:C:231:VAL:HG22	1.84	0.77

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:222:LEU:CD2	1:C:226:GLU:OE2	2.32	0.77
1:B:265:VAL:HG11	1:B:437:THR:HG1	1.50	0.76
1:C:331:TRP:CD2	1:C:336:THR:HG22	2.22	0.75
1:C:106:HIS:CG	1:C:107:PRO:HD2	2.22	0.75
1:C:219:GLY:HA2	1:C:220:PHE:CG	2.20	0.75
1:C:258:TRP:HD1	1:C:258:TRP:H	1.33	0.75
1:B:294:ASN:CB	1:B:384:ALA:O	2.34	0.75
1:B:136:LYS:HD2	1:B:167:ASP:HB3	1.70	0.74
1:C:331:TRP:CG	1:C:336:THR:HG22	2.23	0.74
1:C:258:TRP:CG	1:C:259:TYR:H	2.05	0.73
1:B:132:TYR:CE2	1:B:137:TRP:NE1	2.40	0.73
1:B:263:ILE:HG21	1:B:440:MET:HE1	1.42	0.73
1:B:132:TYR:HE2	1:B:137:TRP:HE1	0.80	0.73
1:B:440:MET:HB3	1:B:441:GLU:CA	2.16	0.73
1:C:68:ARG:HG2	1:C:230:SER:O	1.88	0.73
1:C:258:TRP:H	1:C:258:TRP:CD1	2.06	0.72
1:A:221:PRO:CG	1:A:371:GLU:HG2	2.18	0.72
1:C:144:ASP:OD1	1:C:145:LEU:N	2.22	0.72
1:A:220:PHE:O	1:A:221:PRO:C	2.33	0.71
1:B:279:LYS:CG	1:B:442:ASP:O	2.37	0.71
1:C:332:GLN:HB2	1:C:335:THR:OG1	1.90	0.71
1:A:221:PRO:HG3	1:A:371:GLU:HG2	1.72	0.71
1:C:315:THR:HG23	3:C:615:HOH:O	1.91	0.69
1:B:439:ASP:O	1:B:440:MET:HB2	1.93	0.69
1:B:136:LYS:HB2	1:B:167:ASP:O	1.92	0.69
1:B:278:CYS:SG	1:B:438:LEU:HD23	2.32	0.69
1:B:365:GLN:O	1:B:397:VAL:HB	1.92	0.69
1:B:265:VAL:HG12	1:B:440:MET:CE	2.22	0.69
1:B:350:GLY:H	1:B:356:SER:HB3	1.56	0.68
1:B:353:THR:CG2	1:B:437:THR:O	2.40	0.68
1:C:291:GLY:O	2:F:2:ILE:HG13	1.93	0.68
1:B:363:PRO:HA	1:B:366:TYR:HE2	1.49	0.68
1:B:263:ILE:HG22	1:B:440:MET:HE1	1.75	0.68
1:A:192:ASP:O	1:A:194:LEU:N	2.27	0.68
1:B:263:ILE:HG22	1:B:440:MET:CE	2.21	0.68
1:A:364:GLN:HE22	1:A:424:ASP:HB2	1.58	0.68
1:B:363:PRO:HA	1:B:366:TYR:CD2	2.24	0.67
1:C:417:VAL:CG1	1:C:421:HIS:NE2	2.58	0.67
1:C:219:GLY:HA2	1:C:220:PHE:CD1	2.30	0.67
1:C:67:LEU:HB2	1:C:233:GLY:C	2.21	0.66
1:C:220:PHE:CB	1:C:369:PRO:CB	2.72	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:222:LEU:CD2	1:C:226:GLU:CD	2.67	0.66
1:B:294:ASN:CG	1:B:384:ALA:O	2.39	0.66
1:A:338:TRP:HZ3	1:A:363:PRO:C	2.03	0.66
1:B:130:VAL:HB	1:B:137:TRP:CZ2	2.31	0.66
1:C:377:GLN:HE21	1:C:377:GLN:HA	1.61	0.65
1:A:221:PRO:HB2	1:A:371:GLU:HG3	1.77	0.65
1:C:377:GLN:HA	1:C:377:GLN:NE2	2.09	0.65
1:B:322:PHE:CD1	1:B:329:VAL:HG13	2.31	0.65
1:B:341:PHE:HB3	1:B:366:TYR:OH	1.97	0.65
1:A:338:TRP:HZ3	1:A:364:GLN:CA	2.08	0.65
1:B:136:LYS:CB	1:B:167:ASP:HB3	2.27	0.64
1:B:352:VAL:HG12	1:B:354:ASN:H	1.62	0.64
1:B:265:VAL:CG1	1:B:440:MET:CE	2.76	0.64
1:A:272:GLN:HE22	1:B:427:ARG:HD2	1.61	0.64
1:B:199:ASP:OD1	1:B:408:ARG:NH2	2.31	0.64
1:C:106:HIS:ND1	1:C:107:PRO:HD2	2.14	0.63
1:B:265:VAL:CG1	1:B:437:THR:OG1	2.47	0.62
1:B:279:LYS:NZ	1:B:442:ASP:O	2.24	0.62
1:A:314:SER:HB2	1:B:339:ASN:ND2	2.15	0.62
1:B:265:VAL:CB	1:B:437:THR:OG1	2.47	0.62
1:C:221:PRO:HG2	1:C:370:VAL:C	2.24	0.62
1:B:134:GLN:HG3	1:B:134:GLN:O	2.00	0.62
1:B:263:ILE:CB	1:B:440:MET:HE1	2.30	0.61
1:A:72:GLY:O	2:D:1:GLU:N	2.28	0.61
1:C:220:PHE:CG	1:C:369:PRO:HG2	2.35	0.61
1:C:220:PHE:HB2	1:C:369:PRO:CB	2.31	0.61
1:A:82:GLY:O	1:A:85:PRO:HA	2.00	0.61
1:B:152:PRO:O	1:B:154:VAL:HG22	1.99	0.61
1:B:263:ILE:HG21	1:B:440:MET:SD	2.40	0.60
1:C:219:GLY:CA	1:C:220:PHE:CD1	2.85	0.60
1:A:331:TRP:O	1:A:378:ASP:HB2	2.01	0.60
1:B:338:TRP:HZ3	1:B:367:LEU:HD12	1.65	0.60
1:A:192:ASP:O	1:A:193:SER:C	2.44	0.60
1:C:344:ILE:HB	1:C:361:ILE:HG22	1.82	0.60
1:C:339:ASN:OD1	1:C:340:ILE:N	2.35	0.60
1:B:136:LYS:HB2	1:B:167:ASP:HB3	1.83	0.60
1:C:230:SER:OG	1:C:231:VAL:CG2	2.49	0.60
1:C:192:ASP:O	1:C:193:SER:C	2.44	0.59
1:B:363:PRO:HB3	1:B:366:TYR:CZ	2.37	0.59
1:A:338:TRP:CZ3	1:A:364:GLN:N	2.70	0.59
1:C:220:PHE:HB2	1:C:369:PRO:HB2	1.83	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:338:TRP:HZ3	1:A:364:GLN:N	2.01	0.59
1:B:214:GLN:NE2	1:B:420:CYS:SG	2.75	0.59
1:A:338:TRP:HZ3	1:A:364:GLN:HA	1.56	0.58
1:B:375:THR:O	1:B:375:THR:HG22	2.03	0.58
1:B:291:GLY:C	2:E:2:ILE:HD12	2.27	0.58
1:C:106:HIS:CE1	1:C:107:PRO:HD2	2.39	0.58
1:B:265:VAL:HG21	1:B:437:THR:HB	1.83	0.58
1:C:231:VAL:HB	1:C:232:GLY:C	2.27	0.58
1:A:401:GLY:HA2	1:A:421:HIS:ND1	2.19	0.58
1:A:123:ASP:OD1	1:A:124:LEU:N	2.36	0.57
1:C:372:ASP:CG	1:C:373:VAL:N	2.62	0.57
1:B:363:PRO:CA	1:B:366:TYR:CZ	2.80	0.56
1:C:231:VAL:HB	1:C:233:GLY:H	1.70	0.56
1:A:103:ALA:CB	1:A:162:ALA:HB1	2.36	0.56
1:B:235:MET:HE2	1:B:237:ILE:HG12	1.86	0.56
1:C:258:TRP:CG	1:C:259:TYR:N	2.73	0.56
1:A:63:MET:HG2	1:A:150:HIS:O	2.06	0.55
1:C:134:GLN:OE1	2:F:3:TIH:HD	2.06	0.55
1:B:258:TRP:CD1	1:B:258:TRP:H	2.24	0.55
1:B:210:LEU:HD11	1:B:239:GLY:HA2	1.88	0.55
1:C:71:SER:HB3	1:C:400:GLU:OE1	2.07	0.55
1:A:221:PRO:HB2	1:A:371:GLU:CG	2.36	0.55
1:B:363:PRO:CA	1:B:366:TYR:HE2	2.10	0.54
1:A:83:SER:O	1:A:119:SER:N	2.36	0.54
1:A:332:GLN:HB3	1:A:335:THR:OG1	2.08	0.54
1:B:440:MET:HE2	1:B:443:CYS:HB2	1.88	0.54
1:C:216:CYS:HB2	1:C:231:VAL:HG11	1.89	0.54
1:B:363:PRO:CB	1:B:366:TYR:OH	2.43	0.54
1:A:339:ASN:ND2	1:B:314:SER:HB2	2.23	0.54
1:B:102:GLY:HA2	1:B:163:ILE:HB	1.89	0.54
1:C:128:VAL:HG13	1:C:137:TRP:HZ3	1.73	0.54
1:B:350:GLY:HA3	1:B:355:GLN:HB2	1.90	0.54
1:C:418:SER:HB3	1:C:421:HIS:HD2	1.72	0.53
1:A:201:LEU:O	1:A:205:THR:OG1	2.20	0.53
1:B:258:TRP:CG	1:B:259:TYR:H	2.26	0.53
1:C:70:LYS:HE2	1:C:73:GLN:HG3	1.91	0.53
1:C:332:GLN:O	1:C:333:ALA:C	2.48	0.53
1:A:115:ARG:NH2	1:A:140:GLU:OE2	2.41	0.53
1:B:87:THR:HG22	1:B:111:ARG:HH12	1.74	0.53
1:A:87:THR:HG22	1:A:111:ARG:HH12	1.73	0.52
1:C:67:LEU:N	1:C:233:GLY:O	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:258:TRP:CD1	1:C:258:TRP:N	2.71	0.52
1:C:175:ASN:ND2	1:C:228:LEU:HD23	2.24	0.52
1:A:192:ASP:C	1:A:194:LEU:N	2.64	0.52
1:B:130:VAL:HB	1:B:137:TRP:CE2	2.44	0.52
1:B:194:LEU:O	1:B:195:GLU:C	2.53	0.52
1:B:436:VAL:HG22	1:C:436:VAL:CG2	2.38	0.52
1:C:259:TYR:CE1	2:F:4:LHE:O	2.62	0.52
1:C:332:GLN:O	1:C:335:THR:N	2.35	0.52
1:C:369:PRO:CG	1:C:381:TYR:CD1	2.89	0.52
1:B:205:THR:HG22	1:B:206:HIS:N	2.24	0.51
1:A:175:ASN:HD21	1:A:228:LEU:HD21	1.75	0.51
1:B:136:LYS:HD2	1:B:167:ASP:CB	2.39	0.51
1:C:219:GLY:CA	1:C:220:PHE:CE2	2.74	0.51
1:A:278:CYS:CB	1:A:443:CYS:HG	2.23	0.51
1:C:332:GLN:O	1:C:335:THR:CB	2.58	0.51
1:C:97:SER:OG	1:C:183:ALA:O	2.29	0.51
1:B:363:PRO:CB	1:B:366:TYR:CZ	2.94	0.51
1:B:106:HIS:O	1:B:107:PRO:C	2.54	0.50
1:B:130:VAL:CG2	1:B:137:TRP:CH2	2.94	0.50
1:B:352:VAL:H	1:B:355:GLN:HB2	1.76	0.50
1:B:356:SER:OG	1:B:437:THR:CB	2.59	0.50
1:A:278:CYS:CB	1:A:443:CYS:SG	3.00	0.50
1:B:79:MET:SD	1:B:90:ILE:HG13	2.51	0.50
1:B:153:ASN:O	1:B:154:VAL:HG13	2.11	0.50
1:B:72:GLY:O	2:E:1:GLU:N	2.40	0.50
1:C:58:GLY:O	1:C:59:SER:OG	2.25	0.50
1:C:174:SER:O	1:C:175:ASN:CB	2.58	0.50
1:B:212:SER:HB3	1:B:403:TYR:CE1	2.46	0.50
1:B:439:ASP:OD2	1:C:266:ARG:CD	2.54	0.50
1:A:314:SER:O	1:A:315:THR:C	2.55	0.50
1:C:401:GLY:O	1:C:421:HIS:CD2	2.65	0.50
1:A:278:CYS:HG	1:A:443:CYS:HG	1.45	0.50
1:A:192:ASP:C	1:A:194:LEU:H	2.20	0.49
1:B:169:PHE:CE1	2:E:4:LHE:H33	2.47	0.49
1:B:341:PHE:HB3	1:B:366:TYR:HH	1.76	0.49
1:C:106:HIS:CG	1:C:107:PRO:CD	2.93	0.49
1:B:440:MET:HA	1:B:442:ASP:N	2.22	0.49
1:C:418:SER:O	1:C:421:HIS:CD2	2.65	0.49
1:C:218:ALA:HA	1:C:422:VAL:HG23	1.93	0.49
1:C:372:ASP:OD1	1:C:373:VAL:HG12	2.13	0.49
1:C:418:SER:HB3	1:C:421:HIS:CD2	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:68:ARG:HD2	1:B:228:LEU:HA	1.94	0.49
1:C:97:SER:OG	1:C:183:ALA:C	2.56	0.49
1:C:192:ASP:O	1:C:194:LEU:N	2.45	0.49
1:C:331:TRP:CZ3	1:C:336:THR:HB	2.47	0.49
1:A:221:PRO:CB	1:A:371:GLU:HG2	2.43	0.49
1:C:64:VAL:O	1:C:65:ASP:HB2	2.12	0.49
1:C:67:LEU:CB	1:C:233:GLY:O	2.57	0.49
1:A:258:TRP:CG	1:A:259:TYR:H	2.31	0.48
1:B:295:LEU:HB2	1:B:398:ILE:HD11	1.94	0.48
1:B:364:GLN:HG2	1:B:422:VAL:HG11	1.96	0.48
1:A:318:PHE:CD1	1:A:329:VAL:HG11	2.38	0.48
1:A:339:ASN:HD21	1:B:314:SER:HB2	1.78	0.48
1:A:401:GLY:HA2	1:A:421:HIS:HD1	1.78	0.48
1:C:82:GLY:HA2	1:C:144:ASP:OD2	2.12	0.48
1:C:329:VAL:O	1:C:380:CYS:HA	2.13	0.48
1:C:299:LYS:HG3	1:C:387:GLN:OE1	2.14	0.48
1:B:258:TRP:H	1:B:258:TRP:HD1	1.59	0.48
1:C:251:TYR:CD1	1:C:412:ARG:HG3	2.48	0.48
1:A:216:CYS:N	1:A:232:GLY:O	2.41	0.48
1:B:222:LEU:HD13	1:B:227:VAL:HG22	1.95	0.48
1:A:65:ASP:CG	1:A:231:VAL:HG11	2.39	0.48
1:A:124:LEU:HD12	1:A:141:LEU:HB3	1.95	0.48
1:A:179:ILE:O	1:A:179:ILE:HG23	2.13	0.48
1:A:369:PRO:O	1:A:370:VAL:CG1	2.61	0.48
1:B:348:LEU:O	1:B:356:SER:HB2	2.14	0.48
1:B:90:ILE:HG21	1:B:180:LEU:HB2	1.96	0.47
1:B:250:TRP:O	1:B:414:GLY:HA2	2.14	0.47
1:C:74:GLY:HA3	1:C:91:LEU:HD11	1.96	0.47
1:C:76:TYR:OH	1:C:175:ASN:ND2	2.23	0.47
1:C:106:HIS:CD2	1:C:107:PRO:HD2	2.49	0.47
1:C:191:ASP:O	1:C:192:ASP:C	2.56	0.47
1:A:337:PRO:O	1:A:340:ILE:HG12	2.15	0.47
1:B:436:VAL:O	1:B:436:VAL:HG13	2.13	0.47
1:B:366:TYR:CE1	1:B:367:LEU:CG	2.93	0.47
1:A:330:CYS:HA	1:A:379:ASP:O	2.14	0.47
1:A:260:TYR:HB3	1:A:413:ILE:HD11	1.95	0.47
1:A:332:GLN:HA	1:A:378:ASP:HB2	1.97	0.46
1:B:214:GLN:CD	1:B:420:CYS:SG	2.98	0.46
1:A:223:ASN:OD1	1:A:371:GLU:OE2	2.33	0.46
1:B:123:ASP:OD1	1:B:125:ARG:N	2.33	0.46
1:C:231:VAL:HB	1:C:232:GLY:CA	2.46	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:60:PHE:CZ	1:C:239:GLY:HA3	2.51	0.46
1:A:63:MET:HG2	1:A:151:GLY:HA2	1.98	0.46
1:B:278:CYS:O	1:B:279:LYS:C	2.59	0.46
1:B:314:SER:O	1:B:315:THR:C	2.57	0.46
1:C:364:GLN:OE1	1:C:364:GLN:N	2.41	0.46
1:B:365:GLN:NE2	1:B:401:GLY:HA3	2.31	0.46
1:C:106:HIS:HB3	1:C:109:LEU:HG	1.97	0.46
1:A:84:PRO:HA	1:A:85:PRO:HD3	1.78	0.46
1:A:210:LEU:HD11	1:A:239:GLY:HA2	1.97	0.46
1:A:424:ASP:OD1	1:A:425:GLU:N	2.49	0.46
1:C:363:PRO:HA	1:C:366:TYR:CD2	2.51	0.46
1:A:117:LEU:HD12	1:A:117:LEU:N	2.31	0.45
1:A:331:TRP:HB2	1:A:336:THR:HG22	1.97	0.45
1:B:279:LYS:HA	1:B:443:CYS:O	2.17	0.45
2:E:3:TIH:CD	2:E:3:TIH:N	2.79	0.45
1:A:258:TRP:H	1:A:258:TRP:CD1	2.35	0.45
1:B:354:ASN:ND2	3:B:621:HOH:O	2.50	0.45
1:B:130:VAL:CG2	1:B:137:TRP:CZ2	2.99	0.45
1:C:65:ASP:HA	1:C:232:GLY:O	2.16	0.45
1:A:410:ARG:CD	3:A:612:HOH:O	2.64	0.45
1:C:283:TYR:HA	1:C:284:ASP:HA	1.81	0.45
1:A:336:THR:HG1	1:A:338:TRP:CD1	2.35	0.45
1:C:120:THR:HB	1:C:157:ARG:HH22	1.82	0.45
1:A:115:ARG:HH11	1:A:162:ALA:CB	2.30	0.45
1:A:336:THR:HG21	1:A:381:TYR:CE1	2.52	0.45
1:A:95:GLY:O	2:D:4:LHE:H49	2.17	0.45
1:A:369:PRO:C	1:A:370:VAL:HG13	2.42	0.45
1:C:322:PHE:CD1	1:C:329:VAL:HG23	2.51	0.45
1:C:137:TRP:HB2	1:C:163:ILE:HG23	1.99	0.45
1:C:255:ARG:HD3	1:C:261:GLU:OE2	2.18	0.45
1:B:350:GLY:C	1:B:352:VAL:N	2.75	0.44
1:A:115:ARG:HH11	1:A:162:ALA:HB1	1.82	0.44
1:B:436:VAL:CG2	1:C:436:VAL:HG21	2.40	0.44
1:A:221:PRO:CB	1:A:371:GLU:CG	2.95	0.44
1:A:410:ARG:HD3	3:A:612:HOH:O	2.17	0.44
2:F:1:GLU:OE2	2:F:3:TIH:HE2	2.18	0.44
1:B:216:CYS:HB2	1:B:232:GLY:O	2.18	0.44
1:A:278:CYS:O	1:A:281:TYR:N	2.42	0.44
1:B:375:THR:O	1:B:375:THR:CG2	2.64	0.44
1:C:260:TYR:CE1	1:C:406:PHE:HD2	2.35	0.44
1:A:192:ASP:OD1	1:A:193:SER:N	2.51	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:THR:O	1:A:336:THR:C	2.61	0.44
1:B:205:THR:HG21	1:B:207:VAL:HG23	1.99	0.43
1:C:313:SER:HA	1:C:340:ILE:HG12	2.00	0.43
1:C:411:LYS:HE2	1:C:411:LYS:HB3	1.94	0.43
1:A:372:ASP:OD1	1:A:373:VAL:N	2.52	0.43
1:C:230:SER:HA	1:C:231:VAL:HA	1.72	0.43
1:A:133:THR:HB	2:D:3:TIH:HB2	2.00	0.43
1:B:356:SER:OG	1:B:437:THR:HB	2.18	0.43
1:C:70:LYS:CE	1:C:73:GLN:HG3	2.48	0.43
1:B:293:THR:OG1	2:E:2:ILE:HG13	2.19	0.43
1:B:363:PRO:O	1:B:364:GLN:C	2.60	0.43
1:A:134:GLN:HB3	2:D:3:TIH:HB3	2.00	0.43
1:A:174:SER:O	1:A:175:ASN:CB	2.66	0.43
1:A:355:GLN:O	1:A:440:MET:HE1	2.19	0.43
1:A:97:SER:OG	1:A:184:TYR:O	2.34	0.43
1:A:115:ARG:HD2	1:A:121:TYR:CZ	2.54	0.43
1:A:336:THR:HG21	1:A:381:TYR:HE1	1.83	0.43
1:A:338:TRP:CH2	1:A:364:GLN:HG3	2.54	0.43
1:B:350:GLY:HA3	1:B:355:GLN:CB	2.49	0.43
1:C:329:VAL:HG11	1:C:331:TRP:CZ2	2.54	0.43
1:B:103:ALA:CB	1:B:162:ALA:HB1	2.49	0.42
1:A:183:ALA:HA	1:A:260:TYR:CE2	2.54	0.42
1:A:258:TRP:H	1:A:258:TRP:HD1	1.66	0.42
1:A:428:THR:O	1:A:428:THR:HG23	2.18	0.42
1:C:358:ARG:NH2	1:C:432:GLU:OE2	2.52	0.42
1:C:376:SER:O	1:C:378:ASP:N	2.49	0.42
1:B:366:TYR:CD1	1:B:367:LEU:HG	2.51	0.42
1:C:344:ILE:HB	1:C:361:ILE:CG2	2.49	0.42
1:B:277:ASP:O	1:B:278:CYS:C	2.62	0.42
1:B:118:SER:HB3	1:B:121:TYR:HB2	2.01	0.42
1:B:212:SER:HB2	1:B:236:ILE:HB	2.01	0.42
1:C:218:ALA:HA	1:C:422:VAL:CG2	2.50	0.42
1:C:420:CYS:O	1:C:421:HIS:C	2.61	0.42
1:B:69:GLY:HA2	1:B:227:VAL:HG13	2.01	0.42
1:B:132:TYR:HE2	1:B:137:TRP:CD1	2.30	0.42
1:C:222:LEU:HD23	1:C:226:GLU:OE2	2.18	0.42
1:A:79:MET:SD	1:A:90:ILE:HG13	2.60	0.41
1:B:130:VAL:HG23	1:B:137:TRP:CH2	2.55	0.41
1:B:221:PRO:HG3	1:B:369:PRO:HB2	2.01	0.41
1:B:222:LEU:CD1	1:B:227:VAL:HG22	2.50	0.41
1:B:263:ILE:HG22	1:B:440:MET:HE2	1.93	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:266:ARG:HG2	1:B:267:VAL:N	2.34	0.41
1:C:104:ALA:HB1	1:C:105:PRO:HD2	2.01	0.41
1:A:171:ILE:HD11	2:D:2:ILE:HG22	2.03	0.41
1:A:250:TRP:O	1:A:414:GLY:HA2	2.19	0.41
1:A:249:LEU:HD23	1:A:416:ALA:HB2	2.02	0.41
1:B:258:TRP:CD1	1:B:258:TRP:N	2.89	0.41
1:C:82:GLY:CA	1:C:144:ASP:OD2	2.69	0.41
1:A:314:SER:HB2	1:B:339:ASN:HD22	1.85	0.41
1:B:116:GLN:HG2	1:B:117:LEU:HD12	2.02	0.41
1:B:195:GLU:HA	1:B:196:PRO:HD3	1.93	0.41
1:C:220:PHE:CG	1:C:369:PRO:CG	3.03	0.41
1:B:263:ILE:HB	1:B:440:MET:HE1	2.01	0.41
1:B:338:TRP:CZ3	1:B:367:LEU:HD12	2.52	0.41
1:B:350:GLY:HA3	1:B:356:SER:N	2.36	0.41
1:A:114:GLN:OE1	1:A:117:LEU:HD22	2.21	0.40
1:A:116:GLN:HG2	1:A:117:LEU:HD12	2.01	0.40
1:C:349:MET:HE2	1:C:440:MET:CB	2.51	0.40
1:A:278:CYS:O	1:A:279:LYS:C	2.62	0.40
1:C:258:TRP:HE1	1:C:261:GLU:HB2	1.87	0.40
1:A:366:TYR:HB2	1:A:385:ILE:HG13	2.04	0.40
1:B:78:GLU:O	1:B:149:PRO:HD2	2.22	0.40
1:B:318:PHE:HD1	1:B:329:VAL:HG11	1.84	0.40
1:B:332:GLN:O	1:B:335:THR:OG1	2.29	0.40
1:A:91:LEU:C	1:A:91:LEU:HD23	2.46	0.40
1:A:145:LEU:HD23	1:A:157:ARG:HB2	2.02	0.40
1:B:132:TYR:HB2	1:B:135:GLY:HA3	2.02	0.40
1:B:324:LEU:HD23	1:B:324:LEU:HA	1.98	0.40
1:C:291:GLY:O	2:F:2:ILE:CG1	2.67	0.40
1:C:418:SER:CB	1:C:421:HIS:HD2	2.34	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	387/390 (99%)	381 (98%)	5 (1%)	1 (0%)	36	54
1	B	387/390 (99%)	379 (98%)	7 (2%)	1 (0%)	36	54
1	C	388/390 (100%)	379 (98%)	7 (2%)	2 (0%)	24	42
2	D	1/4 (25%)	1 (100%)	0	0	100	100
2	E	1/4 (25%)	1 (100%)	0	0	100	100
2	F	1/4 (25%)	1 (100%)	0	0	100	100
All	All	1165/1182 (99%)	1142 (98%)	19 (2%)	4 (0%)	36	54

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	221	PRO
1	C	337	PRO
1	A	131	PRO
1	B	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	332/332 (100%)	332 (100%)	0	100	100
1	B	332/332 (100%)	331 (100%)	1 (0%)	86	93
1	C	332/332 (100%)	329 (99%)	3 (1%)	70	84
2	D	2/2 (100%)	2 (100%)	0	100	100
2	E	2/2 (100%)	1 (50%)	1 (50%)	0	0
2	F	2/2 (100%)	2 (100%)	0	100	100
All	All	1002/1002 (100%)	997 (100%)	5 (0%)	81	90

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	437	THR

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Mol	Chain	Res	Type
1	C	258	TRP
1	C	340	ILE
1	C	377	GLN
2	E	2	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	175	ASN
1	A	224	GLN
1	A	272	GLN
1	A	294	ASN
1	A	365	GLN
1	B	214	GLN
1	B	272	GLN
1	C	175	ASN
1	C	294	ASN
1	C	377	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

6 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	LHE	D	4	2	16,16,16	1.03	1 (6%)	15,20,20	1.36	1 (6%)
2	LHE	F	4	2	16,16,16	1.12	1 (6%)	15,20,20	2.08	3 (20%)
2	TIH	D	3	2	9,10,11	0.75	0	6,12,14	1.03	1 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	TIH	F	3	2	9,10,11	1.20	1 (11%)	6,12,14	1.13	1 (16%)
2	TIH	E	3	2	9,10,11	1.44	2 (22%)	6,12,14	1.05	0
2	LHE	E	4	2	16,16,16	0.90	0	15,20,20	1.53	1 (6%)

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	LHE	D	4	2	-	10/20/20/20	-
2	LHE	F	4	2	-	6/20/20/20	-
2	TIH	D	3	2	-	0/5/6/8	0/1/1/1
2	TIH	F	3	2	-	0/5/6/8	0/1/1/1
2	TIH	E	3	2	-	2/5/6/8	0/1/1/1
2	LHE	E	4	2	-	11/20/20/20	-

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	E	3	TIH	CB-CA	-2.60	1.48	1.53
2	E	3	TIH	CG-SD	2.33	1.79	1.70
2	F	3	TIH	CB-CG	2.17	1.57	1.51
2	F	4	LHE	C21-CA2	-2.14	1.49	1.53
2	D	4	LHE	CA-N5	2.01	1.51	1.46

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	4	LHE	C20-N5-CA	-5.34	104.06	114.00
2	E	4	LHE	C20-N5-CA	-3.62	107.27	114.00
2	F	4	LHE	C25-C20-N5	3.34	120.24	111.84
2	D	4	LHE	C20-N5-CA	-3.08	108.26	114.00
2	F	4	LHE	C22-C21-CA2	-3.05	109.20	115.76
2	D	3	TIH	CE1-CE2-SD	-2.31	107.36	113.02
2	F	3	TIH	CE1-CE2-SD	-2.12	107.83	113.02

There are no chirality outliers.

All (29) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	D	4	LHE	C22-C21-CA2-N
2	D	4	LHE	N5-C20-C25-CA2
2	D	4	LHE	N5-C20-C25-O6
2	D	4	LHE	C27-CA-N5-C20
2	D	4	LHE	C-CA-N5-C20
2	E	4	LHE	C22-C21-CA2-C25
2	E	4	LHE	N5-C20-C25-CA2
2	E	4	LHE	N5-C20-C25-O6
2	E	4	LHE	C28-C27-CA-N5
2	F	4	LHE	C-CA-N5-C20
2	E	4	LHE	C28-C27-CA-C
2	E	4	LHE	O-C-CA-N5
2	E	4	LHE	CA-C27-C28-C29
2	F	4	LHE	CA-C27-C28-C29
2	E	4	LHE	OXT-C-CA-N5
2	E	3	TIH	N-CA-CB-CG
2	D	4	LHE	O-C-CA-N5
2	F	4	LHE	CA2-C21-C22-C23
2	D	4	LHE	OXT-C-CA-N5
2	D	4	LHE	CA2-C21-C22-C24
2	E	4	LHE	C25-C20-N5-CA
2	E	4	LHE	OXT-C-CA-C27
2	E	4	LHE	O-C-CA-C27
2	D	4	LHE	C25-C20-N5-CA
2	D	4	LHE	C22-C21-CA2-C25
2	F	4	LHE	C27-CA-N5-C20
2	E	3	TIH	C-CA-CB-CG
2	F	4	LHE	N5-C20-C25-CA2
2	F	4	LHE	O-C-CA-N5

There are no ring outliers.

6 monomers are involved in 8 short contacts:

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

5.5 Carbohydrates [i](#)

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

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	389/390 (99%)	0.05	23 (5%) 28 21	35, 60, 120, 196	0
1	B	389/390 (99%)	0.14	32 (8%) 17 13	40, 59, 110, 190	0
1	C	390/390 (100%)	0.33	37 (9%) 14 11	35, 64, 138, 192	0
2	D	2/4 (50%)	0.57	0 100 100	51, 51, 51, 58	0
2	E	2/4 (50%)	0.20	0 100 100	51, 51, 51, 52	0
2	F	2/4 (50%)	1.30	0 100 100	62, 62, 62, 72	0
All	All	1174/1182 (99%)	0.17	92 (7%) 19 14	35, 61, 127, 196	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	232	GLY	12.9
1	C	231	VAL	8.6
1	B	439	ASP	8.0
1	B	227	VAL	7.1
1	B	440	MET	7.0
1	B	436	VAL	6.3
1	B	366	TYR	6.0
1	C	218	ALA	5.8
1	C	227	VAL	5.7
1	A	220	PHE	5.5
1	C	222	LEU	5.3
1	B	230	SER	5.2
1	A	421	HIS	5.2
1	C	422	VAL	5.1
1	B	442	ASP	4.8
1	A	227	VAL	4.8
1	B	354	ASN	4.7
1	A	422	VAL	4.7
1	B	367	LEU	4.7

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Mol	Chain	Res	Type	RSRZ
1	A	420	CYS	4.6
1	C	229	ALA	4.6
1	A	222	LEU	4.5
1	B	355	GLN	4.4
1	C	355	GLN	4.4
1	C	58	GLY	4.4
1	C	230	SER	4.4
1	B	438	LEU	4.3
1	C	59	SER	4.2
1	A	219	GLY	4.1
1	C	217	GLY	4.0
1	C	219	GLY	4.0
1	B	137	TRP	3.9
1	A	424	ASP	3.9
1	B	437	THR	3.9
1	B	441	GLU	3.9
1	A	231	VAL	3.9
1	C	220	PHE	3.7
1	B	365	GLN	3.5
1	B	220	PHE	3.5
1	C	226	GLU	3.5
1	C	225	SER	3.5
1	C	233	GLY	3.4
1	C	376	SER	3.4
1	C	228	LEU	3.4
1	C	71	SER	3.3
1	B	422	VAL	3.3
1	C	224	GLN	3.3
1	B	70	LYS	3.2
1	A	373	VAL	3.2
1	A	221	PRO	3.2
1	B	373	VAL	3.2
1	B	447	ILE	3.1
1	C	423	HIS	3.1
1	A	230	SER	3.1
1	A	423	HIS	3.0
1	B	150	HIS	3.0
1	C	337	PRO	2.9
1	B	221	PRO	2.9
1	B	136	LYS	2.9
1	B	219	GLY	2.9
1	B	222	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	228	LEU	2.8
1	B	420	CYS	2.7
1	A	218	ALA	2.6
1	C	68	ARG	2.6
1	C	221	PRO	2.6
1	C	369	PRO	2.6
1	B	223	ASN	2.6
1	B	423	HIS	2.5
1	C	373	VAL	2.5
1	C	374	ALA	2.5
1	A	223	ASN	2.5
1	C	421	HIS	2.5
1	A	447	ILE	2.5
1	C	192	ASP	2.4
1	B	135	GLY	2.4
1	C	70	LYS	2.3
1	A	217	GLY	2.3
1	B	229	ALA	2.3
1	C	223	ASN	2.2
1	C	377	GLN	2.2
1	A	59	SER	2.2
1	B	228	LEU	2.2
1	B	231	VAL	2.1
1	C	334	GLY	2.1
1	A	115	ARG	2.1
1	A	229	ALA	2.1
1	C	335	THR	2.1
1	A	225	SER	2.1
1	C	150	HIS	2.0
1	C	266	ARG	2.0
1	A	426	PHE	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. 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
2	LHE	D	4	17/17	0.88	0.15	37,45,70,80	0
2	LHE	E	4	17/17	0.89	0.15	38,52,81,87	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	LHE	F	4	17/17	0.92	0.13	42,46,77,81	0
2	TIH	F	3	10/11	0.94	0.09	47,55,59,69	0
2	TIH	E	3	10/11	0.94	0.10	35,40,44,53	0
2	TIH	D	3	10/11	0.97	0.07	36,44,47,64	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.