



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 12:57 PM UTC

PDB ID : 2J4Y / pdb_00002j4y
Title : Crystal structure of a rhodopsin stabilizing mutant expressed in mammalian cells
Authors : Standfuss, J.; Xie, G.; Edwards, P.C.; Burghammer, M.; Oprian, D.D.; Schertler, G.F.X.
Deposited on : 2006-09-07
Resolution : 3.40 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
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

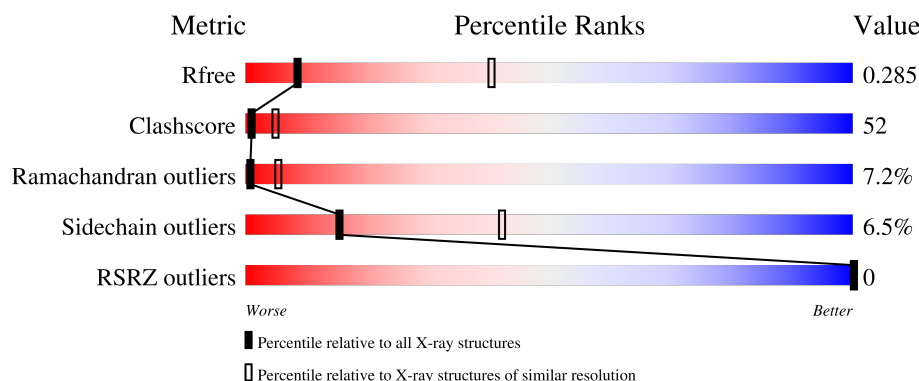
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.40 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	1001 (3.44-3.36)
Clashscore	190562	1022 (3.44-3.36)
Ramachandran outliers	187476	1012 (3.44-3.36)
Sidechain outliers	187428	1012 (3.44-3.36)
RSRZ outliers	180081	1001 (3.44-3.36)

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	349	 28% 52% 12% • 6%
1	B	349	 28% 53% 11% • 6%

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 5252 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.

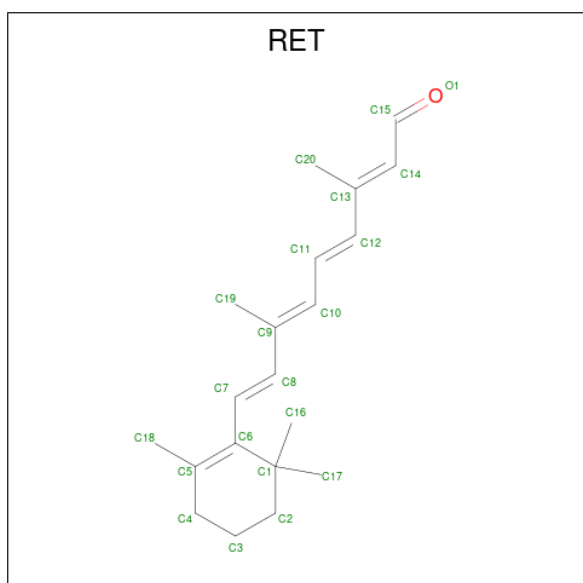
- Molecule 1 is a protein called RHODOPSIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	S	0	0	1
			2592	1725	400	439	28			
1	B	328	Total	C	N	O	S	0	0	1
			2592	1725	400	439	28			

There are 6 discrepancies between the modelled and reference sequences:

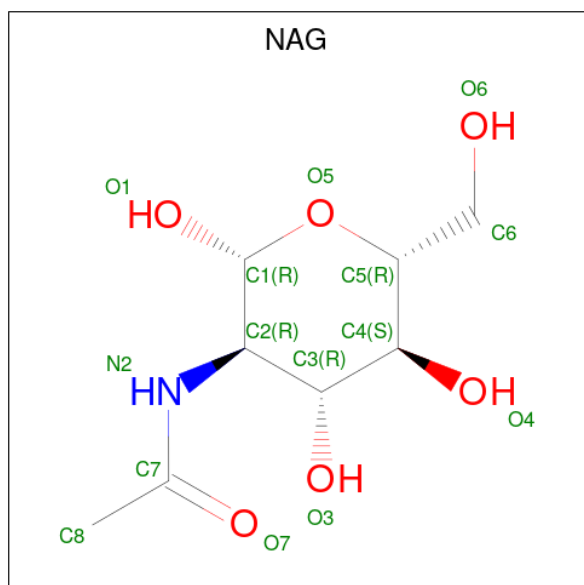
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	ACE	-	acetylation	UNP P02699
A	2	CYS	ASN	engineered mutation	UNP P02699
A	282	CYS	ASP	engineered mutation	UNP P02699
B	0	ACE	-	acetylation	UNP P02699
B	2	CYS	ASN	engineered mutation	UNP P02699
B	282	CYS	ASP	engineered mutation	UNP P02699

- Molecule 2 is RETINAL (CCD ID: RET) (formula: C₂₀H₂₈O).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	C	0	0
			20	20		
2	B	1	Total	C	0	0
			20	20		

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).

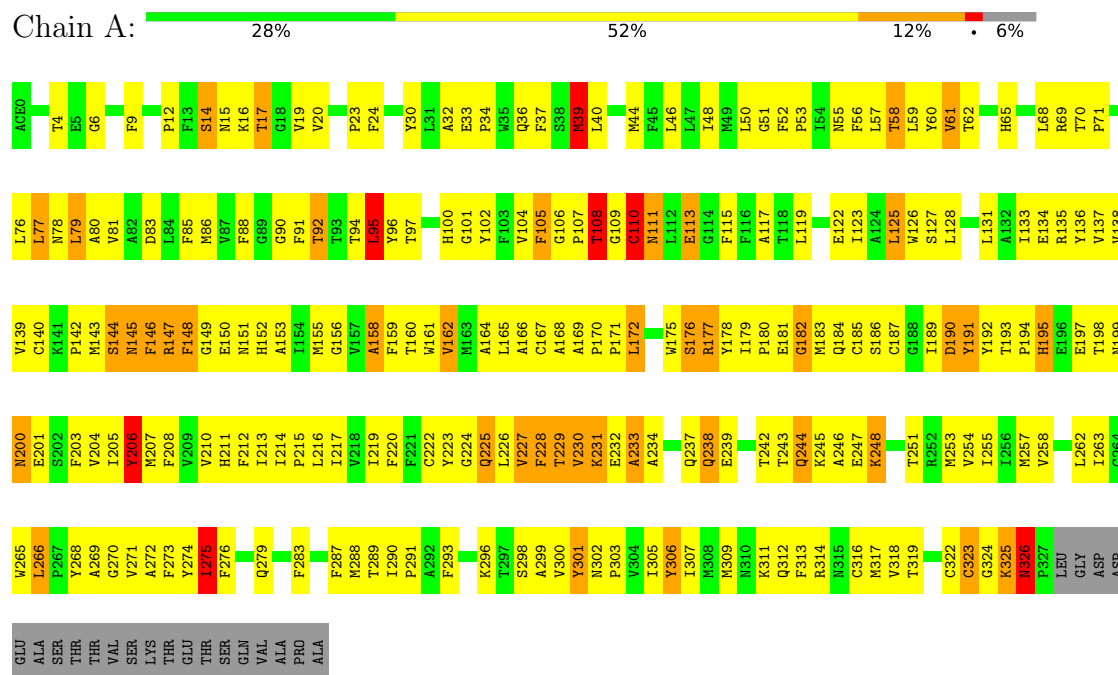


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		

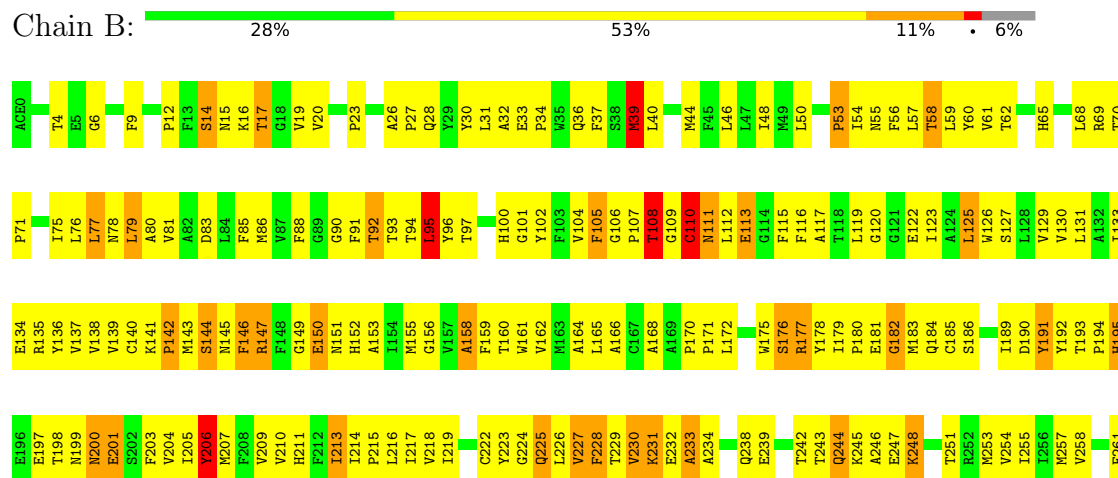
3 Residue-property plots

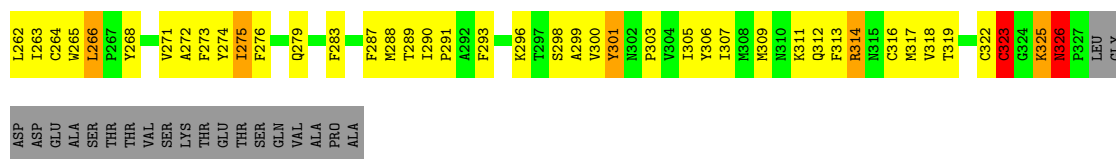
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density.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



• Molecule 1: RHODOPSIN





4 Data and refinement statistics

Property	Value	Source
Space group	P 31	Depositor
Cell constants a, b, c, α , β , γ	109.30Å 109.30Å 77.70Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	50.00 – 3.40 50.00 – 3.40	Depositor EDS
% Data completeness (in resolution range)	95.9 (50.00-3.40) 95.9 (50.00-3.40)	Depositor EDS
R_{merge}	0.24	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.76 (at 3.40Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.290 , 0.330 0.262 , 0.285	Depositor DCC
R_{free} test set	407 reflections (2.97%)	wwPDB-VP
Wilson B-factor (Å ²)	73.3	Xtriage
Anisotropy	0.210	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 80.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.36$, $\langle L^2 \rangle = 0.18$	Xtriage
Estimated twinning fraction	0.340 for -h,-k,l 0.256 for h,-h-k,-l 0.258 for -k,-h,-l	Xtriage
F_o, F_c correlation	0.91	EDS
Total number of atoms	5252	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.47% 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: NAG, RET, ACE

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	0.76	3/2672 (0.1%)	1.20	29/3641 (0.8%)
1	B	0.78	3/2672 (0.1%)	1.19	29/3641 (0.8%)
All	All	0.77	6/5344 (0.1%)	1.20	58/7282 (0.8%)

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

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	15	ASN	CA-CB	6.41	1.60	1.53
1	A	15	ASN	CA-CB	6.32	1.60	1.53
1	B	39	MET	SD-CE	6.32	1.95	1.79
1	B	326	ASN	C-N	-5.61	1.25	1.34
1	A	326	ASN	C-N	-5.47	1.25	1.34
1	A	39	MET	SD-CE	5.23	1.92	1.79

All (58) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	146	PHE	N-CA-C	-9.68	101.69	112.72
1	A	95	LEU	N-CA-C	-8.46	102.01	111.14
1	A	146	PHE	N-CA-C	-8.45	103.46	112.93
1	B	301	TYR	N-CA-C	8.08	121.19	111.82
1	A	15	ASN	N-CA-C	-8.04	102.27	113.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	209	VAL	N-CA-C	7.90	118.70	110.72
1	B	15	ASN	N-CA-C	-7.87	102.50	113.20
1	B	14	SER	N-CA-C	7.77	121.54	109.96
1	B	79	LEU	N-CA-C	-7.76	102.82	111.82
1	B	206	TYR	N-CA-C	-7.76	102.52	110.97
1	A	301	TYR	N-CA-C	7.72	120.78	111.82
1	B	95	LEU	N-CA-C	-7.47	103.07	111.14
1	A	14	SER	N-CA-C	7.42	121.02	109.96
1	A	206	TYR	N-CA-C	-7.35	102.96	110.97
1	A	79	LEU	N-CA-C	-7.17	103.20	112.23
1	A	168	ALA	N-CA-C	7.16	118.87	111.14
1	A	145	ASN	N-CA-C	7.07	120.63	109.39
1	A	117	ALA	N-CA-C	-6.79	104.03	112.72
1	A	246	ALA	N-CA-C	-6.70	104.02	112.93
1	B	117	ALA	N-CA-C	-6.63	104.24	112.72
1	B	145	ASN	N-CA-C	6.58	119.97	108.52
1	A	78	ASN	N-CA-C	-6.41	105.59	113.41
1	A	248	LYS	N-CA-C	-6.37	105.81	113.97
1	B	191	TYR	N-CA-C	-6.34	105.52	113.20
1	B	246	ALA	N-CA-C	-6.22	104.40	112.68
1	B	17	THR	N-CA-C	-6.22	105.48	114.12
1	B	158	ALA	N-CA-C	-6.18	105.59	113.01
1	A	191	TYR	N-CA-C	-6.18	105.72	113.20
1	A	17	THR	N-CA-C	-6.18	105.53	114.12
1	B	265	TRP	N-CA-C	6.17	119.28	111.69
1	B	228	PHE	N-CA-C	-6.09	104.55	111.07
1	A	228	PHE	N-CA-C	-6.04	104.39	110.97
1	B	77	LEU	N-CA-C	-6.02	105.23	113.30
1	A	265	TRP	N-CA-C	5.98	119.04	111.69
1	A	56	PHE	N-CA-C	-5.95	104.37	111.69
1	A	111	ASN	N-CA-C	-5.94	106.07	113.55
1	A	77	LEU	N-CA-C	-5.93	104.72	113.61
1	B	168	ALA	N-CA-C	5.91	118.61	111.40
1	A	158	ALA	N-CA-C	-5.88	105.96	113.01
1	B	78	ASN	N-CA-C	-5.85	106.27	113.41
1	A	325	LYS	CA-CB-CG	-5.85	102.41	114.10
1	B	56	PHE	N-CA-C	-5.79	104.93	112.23
1	A	229	THR	N-CA-C	5.79	117.68	111.36
1	A	251	THR	N-CA-C	-5.78	105.33	112.38
1	A	58	THR	N-CA-C	-5.76	104.98	112.23
1	B	264	CYS	N-CA-C	5.52	117.10	111.14
1	B	251	THR	N-CA-C	-5.47	105.42	111.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	92	THR	N-CA-C	-5.42	105.27	111.07
1	B	111	ASN	N-CA-C	-5.38	106.78	113.55
1	A	275	ILE	CB-CA-C	-5.33	105.15	111.81
1	B	26	ALA	CA-C-N	5.24	125.23	119.89
1	B	26	ALA	C-N-CA	5.24	125.23	119.89
1	A	61	VAL	N-CA-C	-5.20	106.25	113.00
1	A	148	PHE	CB-CA-C	-5.15	103.47	109.80
1	B	58	THR	N-CA-C	-5.13	105.77	112.23
1	B	248	LYS	N-CA-C	-5.07	107.27	113.50
1	B	314	ARG	N-CA-C	5.03	116.84	111.36
1	B	92	THR	N-CA-C	-5.00	105.83	111.28

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	306	TYR	Sidechain

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	2592	0	2567	298	0
1	B	2592	0	2567	284	0
2	A	20	0	27	0	0
2	B	20	0	27	1	0
3	A	14	0	13	0	0
3	B	14	0	13	1	0
All	All	5252	0	5214	547	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 52.

All (547) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:GLY:O	1:A:94:THR:HG23	1.52	1.08
1:A:228:PHE:CE2	1:A:233:ALA:HB2	1.91	1.05
1:B:228:PHE:CE2	1:B:233:ALA:HB2	1.93	1.03
1:A:143:MET:HG2	1:A:144:SER:H	1.22	1.03
1:B:90:GLY:O	1:B:94:THR:HG23	1.58	1.01
1:A:94:THR:HG22	1:A:113:GLU:HG2	1.40	1.00
1:B:194:PRO:O	1:B:195:HIS:O	1.79	1.00
1:A:194:PRO:O	1:A:195:HIS:O	1.79	1.00
1:B:94:THR:HG22	1:B:113:GLU:HG2	1.42	0.97
1:B:143:MET:HG2	1:B:144:SER:H	1.26	0.95
1:A:290:ILE:HB	1:A:291:PRO:HD3	1.47	0.95
1:B:228:PHE:CD2	1:B:233:ALA:HB2	2.02	0.94
1:A:205:ILE:HD13	1:B:228:PHE:HA	1.51	0.93
1:A:228:PHE:HA	1:B:205:ILE:HD13	1.50	0.92
1:B:195:HIS:HB3	1:B:200:ASN:HD22	1.31	0.92
1:A:232:GLU:OE1	1:B:201:GLU:HB2	1.70	0.92
1:B:195:HIS:HB3	1:B:200:ASN:ND2	1.87	0.89
1:A:76:LEU:HD21	1:A:257:MET:HE1	1.52	0.89
1:A:205:ILE:HB	1:B:228:PHE:CD1	2.06	0.89
1:B:290:ILE:HB	1:B:291:PRO:HD3	1.57	0.87
1:B:76:LEU:HD21	1:B:257:MET:HE1	1.56	0.87
1:A:228:PHE:CD1	1:B:205:ILE:HB	2.09	0.86
1:A:319:THR:HG23	1:A:325:LYS:HA	1.58	0.85
1:B:319:THR:HG23	1:B:325:LYS:HA	1.60	0.83
1:A:205:ILE:CB	1:B:228:PHE:HD1	1.91	0.83
1:B:62:THR:HG21	1:B:77:LEU:HD22	1.61	0.83
1:B:143:MET:HG2	1:B:144:SER:N	1.94	0.82
1:A:229:THR:C	1:A:231:LYS:H	1.88	0.81
1:A:230:VAL:HG13	1:A:248:LYS:HE2	1.61	0.81
1:A:228:PHE:HD1	1:B:205:ILE:CB	1.94	0.80
1:B:229:THR:C	1:B:231:LYS:H	1.90	0.80
1:A:228:PHE:CD2	1:A:233:ALA:HB2	2.17	0.78
1:A:86:MET:SD	1:A:299:ALA:HB2	2.24	0.78
1:B:81:VAL:HG12	1:B:85:PHE:CE2	2.20	0.76
1:A:139:VAL:O	1:A:229:THR:HG21	1.86	0.76
1:A:271:VAL:HG13	1:A:287:PHE:CE2	2.21	0.76
1:B:143:MET:CG	1:B:144:SER:H	1.93	0.75
1:A:195:HIS:HB3	1:A:200:ASN:HD22	1.52	0.74
1:B:230:VAL:HG13	1:B:248:LYS:HE2	1.68	0.74
1:B:86:MET:SD	1:B:299:ALA:HB2	2.28	0.74
1:A:37:PHE:CD1	1:A:101:GLY:HA2	2.21	0.74
1:A:232:GLU:C	1:A:234:ALA:H	1.95	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:ILE:HB	1:B:228:PHE:HD1	1.47	0.73
1:A:254:VAL:O	1:A:258:VAL:HG23	1.88	0.73
1:B:254:VAL:O	1:B:258:VAL:HG23	1.89	0.73
1:B:109:GLY:C	1:B:111:ASN:H	1.97	0.73
1:A:62:THR:HG21	1:A:77:LEU:HD22	1.71	0.73
1:A:222:CYS:O	1:A:226:LEU:HG	1.89	0.73
1:A:253:MET:HE1	1:A:306:TYR:CD1	2.25	0.72
1:B:214:ILE:HB	1:B:215:PRO:HD3	1.72	0.72
1:A:95:LEU:C	1:A:95:LEU:HD12	2.15	0.72
1:A:189:ILE:HB	1:A:191:TYR:CE1	2.25	0.72
1:A:214:ILE:HB	1:A:215:PRO:HD3	1.70	0.71
1:B:232:GLU:C	1:B:234:ALA:H	1.98	0.71
1:B:139:VAL:O	1:B:229:THR:HG21	1.90	0.71
1:B:222:CYS:O	1:B:226:LEU:HG	1.90	0.71
1:A:126:TRP:CZ2	1:A:215:PRO:HG3	2.26	0.71
1:A:81:VAL:HG12	1:A:85:PHE:CE2	2.25	0.71
1:A:195:HIS:HB3	1:A:200:ASN:ND2	2.06	0.70
1:A:143:MET:HG2	1:A:144:SER:N	2.01	0.70
1:B:126:TRP:CZ2	1:B:215:PRO:HG3	2.26	0.70
1:A:171:PRO:HD3	1:A:203:PHE:CE2	2.26	0.70
1:A:57:LEU:HD23	1:A:61:VAL:HG23	1.73	0.70
1:B:307:ILE:HG12	1:B:313:PHE:HE2	1.55	0.70
1:A:176:SER:O	1:A:177:ARG:HB3	1.91	0.69
1:A:183:MET:HE3	1:A:289:THR:HG21	1.74	0.69
1:B:95:LEU:HD12	1:B:95:LEU:C	2.16	0.69
1:A:262:LEU:O	1:A:266:LEU:HB2	1.92	0.68
1:B:171:PRO:HD3	1:B:203:PHE:CE2	2.28	0.68
1:B:156:GLY:O	1:B:159:PHE:HB3	1.94	0.68
1:B:271:VAL:HG13	1:B:287:PHE:CE2	2.27	0.68
1:B:262:LEU:O	1:B:266:LEU:HB2	1.92	0.68
1:A:253:MET:HE1	1:A:306:TYR:HD1	1.57	0.68
1:B:228:PHE:CD2	1:B:233:ALA:CB	2.75	0.68
1:B:183:MET:HE3	1:B:289:THR:HG21	1.74	0.68
1:B:189:ILE:HB	1:B:191:TYR:CE1	2.29	0.68
1:A:228:PHE:CD1	1:B:205:ILE:CD1	2.77	0.68
1:A:228:PHE:CD1	1:B:205:ILE:CB	2.74	0.67
1:B:176:SER:O	1:B:177:ARG:HB3	1.93	0.67
1:A:32:ALA:HB1	1:A:36:GLN:OE1	1.95	0.67
1:B:81:VAL:HG12	1:B:85:PHE:HE2	1.57	0.67
1:B:229:THR:HG23	1:B:234:ALA:HB2	1.77	0.67
1:A:109:GLY:C	1:A:111:ASN:H	2.02	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:ILE:HG12	1:A:313:PHE:HE2	1.59	0.67
1:A:205:ILE:CD1	1:B:228:PHE:CD1	2.77	0.67
1:B:225:GLN:C	1:B:227:VAL:H	2.03	0.67
1:A:81:VAL:HG12	1:A:85:PHE:HE2	1.60	0.66
1:A:111:ASN:O	1:A:115:PHE:CB	2.44	0.66
1:B:32:ALA:HB1	1:B:36:GLN:OE1	1.96	0.66
1:A:195:HIS:NE2	1:A:197:GLU:HB2	2.11	0.66
1:A:156:GLY:O	1:A:159:PHE:HB3	1.96	0.66
1:B:194:PRO:HG3	1:B:279:GLN:OE1	1.95	0.66
1:B:262:LEU:HB3	1:B:266:LEU:HD22	1.79	0.65
1:A:262:LEU:HB3	1:A:266:LEU:HD22	1.79	0.65
1:A:318:VAL:HG12	1:A:318:VAL:O	1.96	0.65
1:A:228:PHE:CE1	1:B:205:ILE:HB	2.31	0.65
1:B:150:GLU:O	1:B:153:ALA:HB3	1.96	0.64
1:B:57:LEU:HD23	1:B:61:VAL:HG23	1.79	0.64
1:A:149:GLY:C	1:A:151:ASN:N	2.53	0.64
1:A:225:GLN:C	1:A:227:VAL:H	2.05	0.64
1:B:37:PHE:CD1	1:B:101:GLY:HA2	2.33	0.64
1:B:180:PRO:HB2	1:B:184:GLN:OE1	1.97	0.64
1:A:217:ILE:HG23	1:B:217:ILE:HG23	1.78	0.63
1:A:229:THR:O	1:A:231:LYS:N	2.28	0.63
1:A:205:ILE:HB	1:B:228:PHE:CE1	2.33	0.63
1:B:55:ASN:OD1	1:B:303:PRO:HG2	1.99	0.63
1:B:194:PRO:HA	1:B:201:GLU:OE2	1.98	0.63
1:B:293:PHE:HA	1:B:296:LYS:HE2	1.81	0.63
1:A:76:LEU:HD21	1:A:257:MET:CE	2.26	0.62
1:B:76:LEU:HD21	1:B:257:MET:CE	2.27	0.62
1:A:211:HIS:HA	1:A:215:PRO:HG2	1.80	0.62
1:A:94:THR:CG2	1:A:113:GLU:HG2	2.22	0.62
1:A:324:GLY:C	1:A:325:LYS:HG3	2.25	0.62
1:B:186:SER:HB3	1:B:293:PHE:CE1	2.35	0.62
1:A:134:GLU:O	1:A:138:VAL:HG23	2.00	0.62
1:A:228:PHE:CD2	1:A:233:ALA:CB	2.83	0.62
1:B:94:THR:CG2	1:B:113:GLU:HG2	2.23	0.62
1:B:111:ASN:O	1:B:115:PHE:CB	2.48	0.61
1:A:210:VAL:HG22	1:A:214:ILE:HD12	1.83	0.61
1:A:290:ILE:HB	1:A:291:PRO:CD	2.26	0.61
1:B:253:MET:HE1	1:B:306:TYR:CD1	2.33	0.61
1:A:59:LEU:O	1:A:59:LEU:HG	2.00	0.61
1:A:55:ASN:OD1	1:A:303:PRO:HG2	2.01	0.61
1:A:229:THR:C	1:A:231:LYS:N	2.55	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:134:GLU:O	1:B:138:VAL:HG23	2.00	0.61
1:B:176:SER:H	1:B:198:THR:HB	1.66	0.61
1:B:229:THR:C	1:B:231:LYS:N	2.58	0.61
1:A:217:ILE:CG2	1:B:217:ILE:HG23	2.31	0.61
1:B:20:VAL:HG13	1:B:30:TYR:OH	2.00	0.61
1:B:137:VAL:HG12	1:B:137:VAL:O	1.99	0.60
1:A:115:PHE:CD1	1:A:172:LEU:HD12	2.37	0.60
1:A:303:PRO:O	1:A:307:ILE:HG13	2.02	0.60
1:A:150:GLU:O	1:A:153:ALA:HB3	2.01	0.60
1:B:234:ALA:HB1	1:B:239:GLU:OE1	2.01	0.60
1:B:149:GLY:O	1:B:151:ASN:N	2.35	0.59
1:A:180:PRO:HB2	1:A:184:GLN:OE1	2.02	0.59
1:A:149:GLY:O	1:A:150:GLU:C	2.43	0.59
1:B:107:PRO:C	1:B:109:GLY:H	2.10	0.59
1:A:94:THR:HG21	1:A:113:GLU:OE2	2.02	0.59
1:B:158:ALA:O	1:B:162:VAL:HG23	2.02	0.59
1:B:44:MET:O	1:B:48:ILE:HG13	2.03	0.59
1:A:228:PHE:CE2	1:A:233:ALA:CB	2.79	0.59
1:A:205:ILE:CG2	1:B:228:PHE:HD1	2.16	0.59
1:A:86:MET:HE1	1:A:298:SER:OG	2.03	0.58
1:A:139:VAL:HG12	1:A:229:THR:HB	1.85	0.58
1:A:194:PRO:HG3	1:A:279:GLN:OE1	2.02	0.58
1:A:104:VAL:HG23	1:A:105:PHE:N	2.19	0.58
1:B:318:VAL:HG12	1:B:318:VAL:O	2.03	0.58
1:A:59:LEU:HB2	1:A:80:ALA:HB1	1.86	0.58
1:A:229:THR:HG23	1:A:234:ALA:HB2	1.86	0.58
1:B:17:THR:CG2	1:B:19:VAL:HG12	2.34	0.58
1:A:107:PRO:C	1:A:109:GLY:H	2.12	0.58
1:A:325:LYS:O	1:A:326:ASN:C	2.47	0.58
1:B:14:SER:OG	1:B:16:LYS:HB3	2.04	0.58
1:A:123:ILE:O	1:A:123:ILE:HG22	2.03	0.57
1:A:176:SER:H	1:A:198:THR:HB	1.68	0.57
1:B:123:ILE:HG22	1:B:123:ILE:O	2.04	0.57
1:B:149:GLY:C	1:B:151:ASN:N	2.59	0.57
1:A:17:THR:CG2	1:A:19:VAL:HG12	2.34	0.57
1:A:232:GLU:CD	1:B:201:GLU:HB2	2.29	0.57
1:B:211:HIS:HA	1:B:215:PRO:HG2	1.86	0.57
1:B:307:ILE:O	1:B:314:ARG:HD3	2.04	0.57
1:B:83:ASP:O	1:B:86:MET:HB2	2.05	0.57
1:B:210:VAL:HG22	1:B:214:ILE:HD12	1.85	0.57
1:A:149:GLY:O	1:A:151:ASN:N	2.37	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:234:ALA:HB1	1:A:239:GLU:OE1	2.04	0.57
1:B:62:THR:CG2	1:B:77:LEU:HD22	2.33	0.57
1:A:17:THR:HG23	1:A:19:VAL:HG12	1.86	0.57
1:A:271:VAL:HG12	1:A:288:MET:HE2	1.86	0.57
1:A:210:VAL:O	1:A:215:PRO:HD3	2.03	0.57
1:B:253:MET:HE1	1:B:306:TYR:HD1	1.70	0.57
1:A:111:ASN:O	1:A:115:PHE:HB2	2.05	0.56
1:A:217:ILE:HG23	1:B:217:ILE:CG2	2.34	0.56
1:A:228:PHE:HD1	1:B:205:ILE:CD1	2.17	0.56
1:A:228:PHE:HD1	1:B:205:ILE:CG2	2.18	0.56
1:A:57:LEU:HD23	1:A:61:VAL:CG2	2.35	0.56
1:A:166:ALA:O	1:A:170:PRO:HG2	2.06	0.56
1:A:179:ILE:HG21	1:A:190:ASP:OD1	2.06	0.56
1:A:276:PHE:O	1:A:279:GLN:HG3	2.05	0.56
1:A:317:MET:HE2	1:A:318:VAL:HG23	1.85	0.56
1:B:176:SER:O	1:B:177:ARG:CB	2.54	0.56
1:B:232:GLU:C	1:B:234:ALA:N	2.63	0.56
1:B:307:ILE:HG12	1:B:313:PHE:CE2	2.38	0.56
1:B:317:MET:HE2	1:B:318:VAL:HG23	1.87	0.56
1:A:135:ARG:HB3	1:A:226:LEU:HD21	1.87	0.56
1:B:104:VAL:HG23	1:B:105:PHE:N	2.21	0.56
1:A:268:TYR:HA	1:A:291:PRO:HB3	1.87	0.56
1:B:192:TYR:CE1	1:B:275:ILE:HD13	2.40	0.56
1:A:65:HIS:CE1	1:A:316:CYS:HB3	2.41	0.56
1:B:106:GLY:C	1:B:108:THR:N	2.63	0.56
1:B:175:TRP:O	1:B:176:SER:HB2	2.06	0.56
1:A:76:LEU:HD23	1:A:79:LEU:HD12	1.88	0.56
1:A:307:ILE:O	1:A:314:ARG:HD3	2.05	0.56
1:A:137:VAL:HG12	1:A:137:VAL:O	2.05	0.55
1:B:23:PRO:HB3	1:B:184:GLN:HB3	1.88	0.55
1:B:37:PHE:HE1	1:B:185:CYS:SG	2.29	0.55
1:A:111:ASN:O	1:A:115:PHE:HB3	2.04	0.55
1:A:203:PHE:CD1	1:A:203:PHE:C	2.84	0.55
1:B:322:CYS:O	1:B:323:CYS:O	2.25	0.55
1:A:44:MET:O	1:A:48:ILE:HG13	2.07	0.55
1:B:276:PHE:O	1:B:279:GLN:HG3	2.06	0.55
1:B:94:THR:HG21	1:B:113:GLU:OE2	2.07	0.55
1:B:143:MET:CG	1:B:144:SER:N	2.62	0.55
1:A:186:SER:HB3	1:A:293:PHE:CE1	2.42	0.55
1:A:83:ASP:O	1:A:86:MET:HB2	2.07	0.54
1:A:205:ILE:CB	1:B:228:PHE:CD1	2.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:14:SER:OG	1:A:16:LYS:HB3	2.07	0.54
1:A:318:VAL:O	1:A:318:VAL:CG1	2.54	0.54
1:B:76:LEU:HD23	1:B:79:LEU:HD12	1.88	0.54
1:B:149:GLY:O	1:B:150:GLU:C	2.50	0.54
1:B:166:ALA:O	1:B:170:PRO:HG2	2.07	0.54
1:A:205:ILE:CD1	1:B:228:PHE:HD1	2.19	0.54
1:A:207:MET:O	1:A:211:HIS:HB2	2.07	0.54
1:B:12:PRO:HB2	1:B:283:PHE:O	2.08	0.54
1:B:146:PHE:O	1:B:147:ARG:C	2.49	0.54
1:A:88:PHE:HA	1:A:92:THR:HG23	1.89	0.54
1:A:232:GLU:C	1:A:234:ALA:N	2.62	0.54
1:A:293:PHE:HA	1:A:296:LYS:HE2	1.89	0.54
1:B:195:HIS:NE2	1:B:197:GLU:HB2	2.22	0.54
1:B:231:LYS:HG3	1:B:232:GLU:N	2.23	0.54
1:A:110:CYS:SG	1:A:110:CYS:O	2.65	0.54
1:B:91:PHE:CZ	1:B:296:LYS:HG2	2.42	0.54
1:B:111:ASN:O	1:B:115:PHE:HB3	2.08	0.54
1:A:12:PRO:HB2	1:A:283:PHE:O	2.08	0.53
1:A:192:TYR:OH	1:A:288:MET:HG2	2.08	0.53
1:A:243:THR:HG22	1:A:247:GLU:HG3	1.89	0.53
1:B:69:ARG:HG2	1:B:69:ARG:HH11	1.73	0.53
1:B:192:TYR:OH	1:B:288:MET:HG2	2.09	0.53
1:B:290:ILE:HB	1:B:291:PRO:CD	2.34	0.53
1:A:106:GLY:C	1:A:108:THR:N	2.64	0.53
1:A:231:LYS:HG3	1:A:232:GLU:N	2.23	0.53
1:A:271:VAL:HG11	1:A:287:PHE:O	2.08	0.53
1:B:152:HIS:HA	1:B:155:MET:CG	2.38	0.53
1:B:59:LEU:O	1:B:59:LEU:HG	2.09	0.53
1:B:88:PHE:HA	1:B:92:THR:HG23	1.90	0.53
1:A:77:LEU:HG	1:A:77:LEU:O	2.08	0.53
1:A:176:SER:O	1:A:177:ARG:CB	2.53	0.53
1:A:192:TYR:CE1	1:A:275:ILE:HD13	2.44	0.53
1:A:317:MET:O	1:A:317:MET:HG2	2.09	0.53
1:B:119:LEU:CD1	1:B:165:LEU:HD23	2.39	0.53
1:B:127:SER:OG	1:B:160:THR:HG21	2.09	0.53
1:A:158:ALA:O	1:A:162:VAL:HG23	2.09	0.53
1:B:57:LEU:HD23	1:B:61:VAL:CG2	2.39	0.53
1:B:130:VAL:HG12	1:B:130:VAL:O	2.09	0.53
1:B:216:LEU:O	1:B:219:ILE:N	2.42	0.53
1:A:127:SER:OG	1:A:160:THR:HG21	2.09	0.52
1:A:205:ILE:HD13	1:B:228:PHE:CD1	2.44	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:307:ILE:HG12	1:A:313:PHE:CE2	2.43	0.52
1:B:135:ARG:HB3	1:B:226:LEU:HD21	1.91	0.52
1:A:20:VAL:HG13	1:A:30:TYR:OH	2.08	0.52
1:A:205:ILE:HD12	1:B:228:PHE:CD1	2.45	0.52
1:A:228:PHE:CD1	1:B:205:ILE:HG21	2.44	0.52
1:B:70:THR:HB	1:B:71:PRO:HD2	1.91	0.52
1:B:75:ILE:HG21	1:B:131:LEU:HD22	1.91	0.52
1:B:106:GLY:O	1:B:108:THR:N	2.42	0.52
1:A:109:GLY:C	1:A:111:ASN:N	2.68	0.52
1:B:223:TYR:O	1:B:226:LEU:HB2	2.09	0.52
1:A:70:THR:HB	1:A:71:PRO:HD2	1.90	0.52
1:B:20:VAL:HA	1:B:30:TYR:CZ	2.45	0.52
1:A:119:LEU:CD1	1:A:165:LEU:HD23	2.40	0.52
1:A:223:TYR:O	1:A:226:LEU:HB2	2.10	0.52
1:B:271:VAL:HG12	1:B:288:MET:HE2	1.91	0.52
1:A:91:PHE:CZ	1:A:296:LYS:HG2	2.44	0.52
1:B:57:LEU:O	1:B:61:VAL:HG23	2.10	0.52
1:B:109:GLY:C	1:B:111:ASN:N	2.64	0.52
1:B:133:ILE:O	1:B:137:VAL:HG23	2.10	0.52
1:B:268:TYR:HA	1:B:291:PRO:HB3	1.92	0.52
1:A:272:ALA:N	1:A:288:MET:HE1	2.26	0.51
1:B:181:GLU:HB2	1:B:186:SER:O	2.09	0.51
1:B:325:LYS:O	1:B:326:ASN:C	2.51	0.51
1:A:205:ILE:CG2	1:B:228:PHE:CD1	2.93	0.51
1:A:228:PHE:CD1	1:B:205:ILE:CG2	2.93	0.51
1:A:228:PHE:CD1	1:B:205:ILE:HD12	2.45	0.51
1:B:171:PRO:HB3	1:B:176:SER:HB3	1.92	0.51
1:B:311:LYS:O	1:B:312:GLN:C	2.53	0.51
1:B:192:TYR:CE1	1:B:275:ILE:CD1	2.94	0.51
1:B:229:THR:O	1:B:231:LYS:N	2.37	0.51
1:A:205:ILE:HG21	1:B:228:PHE:CD1	2.46	0.51
1:A:59:LEU:O	1:A:59:LEU:CG	2.59	0.51
1:A:139:VAL:HG21	1:A:226:LEU:HD22	1.93	0.51
1:B:152:HIS:HA	1:B:155:MET:HG2	1.93	0.51
1:B:161:TRP:O	1:B:165:LEU:HG	2.10	0.51
1:B:200:ASN:O	1:B:204:VAL:HG23	2.10	0.51
1:A:24:PHE:CE1	1:A:107:PRO:HA	2.46	0.51
1:B:243:THR:C	1:B:245:LYS:H	2.19	0.51
1:A:135:ARG:HB3	1:A:226:LEU:CD2	2.40	0.51
1:A:243:THR:O	1:A:247:GLU:HB2	2.11	0.51
1:B:33:GLU:OE1	1:B:34:PRO:HD2	2.11	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:145:ASN:C	1:A:147:ARG:H	2.18	0.51
1:B:215:PRO:O	1:B:219:ILE:HG13	2.11	0.51
1:B:318:VAL:O	1:B:318:VAL:CG1	2.59	0.51
1:A:322:CYS:O	1:A:323:CYS:O	2.28	0.50
1:B:303:PRO:O	1:B:307:ILE:HG13	2.11	0.50
1:A:228:PHE:CD1	1:B:205:ILE:HD13	2.45	0.50
1:A:243:THR:CG2	1:A:247:GLU:HG3	2.40	0.50
1:B:106:GLY:C	1:B:108:THR:H	2.18	0.50
1:B:139:VAL:HG12	1:B:229:THR:HB	1.93	0.50
1:B:203:PHE:C	1:B:203:PHE:CD1	2.90	0.50
1:A:133:ILE:O	1:A:137:VAL:HG23	2.12	0.50
1:B:59:LEU:HB2	1:B:80:ALA:HB1	1.94	0.50
1:A:70:THR:HB	1:A:71:PRO:CD	2.41	0.50
1:A:135:ARG:HA	1:A:135:ARG:NE	2.26	0.50
1:A:178:TYR:CE1	1:A:189:ILE:HD13	2.46	0.50
1:B:179:ILE:HG21	1:B:190:ASP:OD1	2.10	0.50
1:A:146:PHE:O	1:A:147:ARG:C	2.54	0.50
1:B:224:GLY:O	1:B:227:VAL:N	2.44	0.50
1:B:224:GLY:O	1:B:225:GLN:C	2.55	0.50
1:A:4:THR:HB	1:A:20:VAL:HG11	1.94	0.49
1:A:244:GLN:OE1	1:A:248:LYS:HE3	2.12	0.49
1:B:111:ASN:O	1:B:115:PHE:HB2	2.11	0.49
1:B:40:LEU:HD22	1:B:293:PHE:CE2	2.48	0.49
1:A:106:GLY:O	1:A:108:THR:N	2.45	0.49
1:A:181:GLU:HB2	1:A:186:SER:O	2.12	0.49
1:B:48:ILE:CD1	1:B:95:LEU:HD23	2.43	0.49
1:A:62:THR:CG2	1:A:77:LEU:HD22	2.40	0.49
1:B:37:PHE:CE1	1:B:185:CYS:SG	3.06	0.49
1:A:143:MET:CG	1:A:144:SER:H	1.95	0.49
1:B:40:LEU:HD22	1:B:293:PHE:CD2	2.48	0.49
1:B:243:THR:HG22	1:B:247:GLU:HG3	1.93	0.49
1:B:272:ALA:N	1:B:288:MET:HE1	2.27	0.49
1:A:152:HIS:HA	1:A:155:MET:CG	2.42	0.49
1:B:176:SER:HB2	1:B:203:PHE:CG	2.48	0.49
1:A:228:PHE:HD1	1:B:205:ILE:HG21	1.77	0.49
1:B:135:ARG:HB3	1:B:226:LEU:CD2	2.43	0.49
1:A:106:GLY:C	1:A:108:THR:H	2.21	0.49
1:A:210:VAL:HA	1:A:214:ILE:HD12	1.95	0.49
1:B:57:LEU:HD23	1:B:57:LEU:O	2.13	0.49
1:A:44:MET:HE2	1:A:94:THR:OG1	2.13	0.48
1:A:242:THR:O	1:A:242:THR:HG22	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:224:GLY:O	1:A:226:LEU:N	2.46	0.48
1:A:305:ILE:HG23	1:A:309:MET:SD	2.52	0.48
1:B:136:TYR:HE1	1:B:225:GLN:HE21	1.61	0.48
1:B:170:PRO:HB2	1:B:203:PHE:CE2	2.48	0.48
1:B:307:ILE:HD13	1:B:317:MET:SD	2.52	0.48
1:A:149:GLY:C	1:A:151:ASN:H	2.20	0.48
1:B:139:VAL:HG21	1:B:226:LEU:HD22	1.95	0.48
1:B:48:ILE:HD12	1:B:95:LEU:HD23	1.95	0.48
1:A:106:GLY:O	1:A:107:PRO:C	2.56	0.48
1:A:215:PRO:O	1:A:219:ILE:HG13	2.13	0.48
1:B:176:SER:HB2	1:B:203:PHE:CD1	2.49	0.48
1:A:136:TYR:HE1	1:A:225:GLN:HE21	1.61	0.48
1:B:106:GLY:O	1:B:107:PRO:C	2.55	0.48
1:A:194:PRO:O	1:A:195:HIS:C	2.55	0.48
1:B:17:THR:HG23	1:B:19:VAL:HG12	1.95	0.48
1:B:149:GLY:C	1:B:151:ASN:H	2.20	0.48
1:A:208:PHE:O	1:A:212:PHE:HB3	2.14	0.48
1:A:14:SER:C	1:A:16:LYS:H	2.22	0.47
1:A:135:ARG:O	1:A:139:VAL:HG23	2.14	0.47
1:A:139:VAL:HG21	1:A:226:LEU:CD2	2.44	0.47
1:B:94:THR:HG22	1:B:113:GLU:CG	2.30	0.47
1:B:125:LEU:HD22	1:B:215:PRO:HB3	1.96	0.47
1:A:6:GLY:HA3	1:A:9:PHE:CZ	2.49	0.47
1:A:44:MET:HE1	1:A:91:PHE:CD1	2.49	0.47
1:A:224:GLY:O	1:A:227:VAL:N	2.46	0.47
1:B:14:SER:C	1:B:16:LYS:H	2.22	0.47
1:B:224:GLY:O	1:B:226:LEU:N	2.47	0.47
1:A:171:PRO:HB3	1:A:176:SER:HB3	1.96	0.47
1:A:9:PHE:C	1:A:179:ILE:HD11	2.40	0.47
1:A:91:PHE:CE1	1:A:296:LYS:HD3	2.50	0.47
1:A:179:ILE:C	1:A:187:CYS:SG	2.98	0.47
1:B:81:VAL:CG1	1:B:85:PHE:HE2	2.27	0.47
1:A:224:GLY:O	1:A:225:GLN:C	2.56	0.47
1:A:170:PRO:HB2	1:A:203:PHE:CE2	2.50	0.47
1:A:175:TRP:O	1:A:176:SER:HB2	2.14	0.47
1:A:201:GLU:HG2	1:B:231:LYS:HE3	1.97	0.47
1:A:225:GLN:C	1:A:227:VAL:N	2.73	0.47
1:A:272:ALA:CA	1:A:288:MET:HE1	2.45	0.47
1:A:131:LEU:HD13	1:A:254:VAL:HG22	1.96	0.47
1:A:166:ALA:O	1:A:206:TYR:OH	2.32	0.47
1:B:210:VAL:HA	1:B:214:ILE:HD12	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:LEU:HD13	1:A:306:TYR:CD1	2.50	0.47
1:B:271:VAL:O	1:B:272:ALA:C	2.58	0.47
1:A:205:ILE:HG21	1:B:228:PHE:HB2	1.96	0.46
1:A:216:LEU:O	1:A:219:ILE:N	2.47	0.46
1:B:179:ILE:HG12	1:B:180:PRO:N	2.30	0.46
1:A:20:VAL:HA	1:A:30:TYR:CZ	2.50	0.46
1:A:76:LEU:HD11	1:A:253:MET:HE2	1.97	0.46
1:B:77:LEU:O	1:B:77:LEU:HG	2.15	0.46
1:A:19:VAL:HG13	1:A:20:VAL:N	2.31	0.46
1:A:217:ILE:CG2	1:B:217:ILE:CG2	2.92	0.46
1:A:271:VAL:O	1:A:272:ALA:C	2.58	0.46
1:A:311:LYS:O	1:A:312:GLN:C	2.58	0.46
1:A:37:PHE:CE1	1:A:101:GLY:HA2	2.51	0.46
1:A:113:GLU:C	1:A:113:GLU:CD	2.83	0.46
1:B:91:PHE:HA	1:B:94:THR:OG1	2.16	0.46
1:B:129:VAL:HG13	1:B:218:VAL:CG1	2.44	0.46
1:B:135:ARG:NE	1:B:135:ARG:HA	2.29	0.46
1:A:169:ALA:O	1:A:170:PRO:C	2.59	0.46
1:B:6:GLY:HA3	1:B:9:PHE:CZ	2.50	0.46
1:B:86:MET:HE1	1:B:298:SER:OG	2.16	0.46
1:B:110:CYS:O	1:B:110:CYS:SG	2.73	0.46
1:A:181:GLU:O	1:A:182:GLY:C	2.58	0.46
1:B:59:LEU:O	1:B:59:LEU:CG	2.64	0.46
1:B:181:GLU:O	1:B:182:GLY:C	2.59	0.46
1:A:179:ILE:HG12	1:A:180:PRO:N	2.29	0.46
1:A:301:TYR:CD1	1:A:301:TYR:C	2.93	0.46
1:B:271:VAL:HG11	1:B:287:PHE:O	2.16	0.46
1:B:9:PHE:C	1:B:179:ILE:HD11	2.40	0.46
1:A:176:SER:HB2	1:A:203:PHE:CD1	2.51	0.46
1:B:170:PRO:HB2	1:B:203:PHE:HE2	1.81	0.46
1:B:207:MET:O	1:B:211:HIS:HB2	2.16	0.46
1:A:79:LEU:O	1:A:83:ASP:CG	2.59	0.46
1:A:192:TYR:CE1	1:A:275:ILE:CD1	2.98	0.46
1:B:170:PRO:N	1:B:171:PRO:HD2	2.30	0.46
1:A:33:GLU:OE1	1:A:34:PRO:HD2	2.15	0.45
1:A:57:LEU:O	1:A:61:VAL:HG23	2.16	0.45
1:A:69:ARG:HG2	1:A:69:ARG:HH11	1.81	0.45
1:A:170:PRO:N	1:A:171:PRO:HD2	2.31	0.45
1:B:232:GLU:O	1:B:234:ALA:N	2.48	0.45
1:A:140:CYS:HB3	1:A:229:THR:OG1	2.15	0.45
1:A:152:HIS:HA	1:A:155:MET:HG2	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:70:THR:HB	1:B:71:PRO:CD	2.45	0.45
1:A:122:GLU:HB3	1:A:164:ALA:HA	1.97	0.45
1:A:135:ARG:CB	1:A:226:LEU:HD21	2.47	0.45
1:B:180:PRO:HB3	1:B:184:GLN:HA	1.98	0.45
1:A:125:LEU:HD22	1:A:215:PRO:HB3	1.98	0.45
1:A:128:LEU:HD23	1:A:128:LEU:HA	1.78	0.45
1:B:190:ASP:C	1:B:192:TYR:H	2.25	0.45
1:B:175:TRP:O	1:B:176:SER:CB	2.64	0.45
1:B:243:THR:CG2	1:B:247:GLU:HG3	2.46	0.45
3:B:1327:NAG:O7	3:B:1327:NAG:O3	2.28	0.45
1:A:232:GLU:O	1:A:232:GLU:HG2	2.17	0.45
1:B:76:LEU:HD11	1:B:253:MET:HE2	1.98	0.45
1:A:76:LEU:CD2	1:A:257:MET:HE1	2.37	0.45
1:B:242:THR:HG22	1:B:242:THR:O	2.15	0.45
1:B:91:PHE:CE1	1:B:296:LYS:HG2	2.52	0.44
1:B:199:ASN:O	1:B:200:ASN:C	2.59	0.44
1:A:176:SER:HB2	1:A:203:PHE:CG	2.52	0.44
1:A:199:ASN:O	1:A:200:ASN:C	2.59	0.44
1:B:4:THR:HB	1:B:20:VAL:HG11	1.98	0.44
1:B:19:VAL:HG13	1:B:20:VAL:N	2.33	0.44
1:B:76:LEU:HD13	1:B:306:TYR:CD1	2.52	0.44
1:A:190:ASP:C	1:A:192:TYR:H	2.25	0.44
1:B:178:TYR:CE1	1:B:189:ILE:HD13	2.52	0.44
1:B:243:THR:O	1:B:247:GLU:HB2	2.17	0.44
1:B:86:MET:SD	1:B:120:GLY:HA3	2.58	0.44
1:B:301:TYR:CD1	1:B:301:TYR:C	2.95	0.44
1:A:30:TYR:CD1	1:A:30:TYR:C	2.95	0.44
1:A:37:PHE:HE1	1:A:185:CYS:SG	2.41	0.44
1:B:139:VAL:HG21	1:B:226:LEU:CD2	2.47	0.44
1:B:199:ASN:O	1:B:201:GLU:N	2.51	0.44
1:B:191:TYR:OH	2:B:1296:RET:H192	2.18	0.44
1:A:319:THR:HG22	1:A:324:GLY:O	2.18	0.44
1:B:147:ARG:HH11	1:B:147:ARG:HG2	1.83	0.44
1:A:46:LEU:HD23	1:A:46:LEU:HA	1.88	0.44
1:A:52:PHE:CD1	1:A:88:PHE:HZ	2.36	0.44
1:A:205:ILE:HG21	1:B:228:PHE:HD1	1.78	0.44
1:A:161:TRP:O	1:A:165:LEU:HG	2.18	0.43
1:B:273:PHE:O	1:B:274:TYR:C	2.61	0.43
1:B:79:LEU:HD23	1:B:79:LEU:HA	1.85	0.43
1:B:88:PHE:HA	1:B:92:THR:CG2	2.48	0.43
1:B:225:GLN:C	1:B:227:VAL:N	2.71	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:112:LEU:O	1:B:116:PHE:HB2	2.18	0.43
1:A:95:LEU:HG	1:A:96:TYR:N	2.34	0.43
1:B:113:GLU:CD	1:B:113:GLU:C	2.86	0.43
1:B:244:GLN:OE1	1:B:248:LYS:HE3	2.18	0.43
1:A:51:GLY:O	1:A:55:ASN:HB2	2.19	0.43
1:A:96:TYR:HE2	1:A:104:VAL:HG21	1.83	0.43
1:A:100:HIS:O	1:A:102:TYR:HD1	2.02	0.43
1:A:179:ILE:HG13	1:A:180:PRO:HD2	2.01	0.43
1:A:223:TYR:O	1:A:224:GLY:C	2.62	0.43
1:A:319:THR:CG2	1:A:324:GLY:O	2.67	0.43
1:A:94:THR:HG22	1:A:113:GLU:CG	2.29	0.43
1:B:100:HIS:O	1:B:102:TYR:HD1	2.02	0.43
1:A:39:MET:HE3	1:A:39:MET:HA	1.99	0.43
1:B:68:LEU:HD23	1:B:312:GLN:HE21	1.84	0.43
1:B:141:LYS:HA	1:B:142:PRO:HD3	1.82	0.43
1:A:62:THR:HB	1:A:77:LEU:HD13	2.00	0.42
1:A:81:VAL:CG1	1:A:85:PHE:HE2	2.30	0.42
1:A:88:PHE:HA	1:A:92:THR:CG2	2.49	0.42
1:A:108:THR:O	1:A:108:THR:OG1	2.29	0.42
1:A:175:TRP:O	1:A:176:SER:CB	2.67	0.42
1:A:194:PRO:C	1:A:195:HIS:O	2.56	0.42
1:B:272:ALA:CA	1:B:288:MET:HE1	2.49	0.42
1:A:210:VAL:O	1:A:215:PRO:CD	2.67	0.42
1:B:135:ARG:CB	1:B:226:LEU:HD21	2.48	0.42
1:B:58:THR:C	1:B:60:TYR:H	2.27	0.42
1:B:93:THR:O	1:B:93:THR:HG22	2.20	0.42
1:B:213:ILE:O	1:B:214:ILE:C	2.61	0.42
1:A:232:GLU:O	1:A:234:ALA:N	2.53	0.42
1:B:44:MET:HE1	1:B:91:PHE:CD1	2.55	0.42
1:A:172:LEU:HD23	1:A:172:LEU:O	2.20	0.42
1:A:23:PRO:HB3	1:A:184:GLN:HB3	2.02	0.42
1:A:200:ASN:O	1:A:204:VAL:HG23	2.19	0.42
1:A:266:LEU:HD12	1:A:266:LEU:HA	1.84	0.42
1:B:28:GLN:HB3	1:B:31:LEU:HD12	2.01	0.42
1:B:69:ARG:HG2	1:B:69:ARG:NH1	2.34	0.42
1:A:58:THR:C	1:A:60:TYR:H	2.28	0.42
1:A:257:MET:HE3	1:A:302:ASN:OD1	2.19	0.42
1:B:39:MET:HA	1:B:39:MET:HE3	2.02	0.42
1:B:135:ARG:O	1:B:139:VAL:HG23	2.20	0.42
1:A:91:PHE:HA	1:A:94:THR:OG1	2.20	0.42
1:B:125:LEU:HG	1:B:261:PHE:CZ	2.55	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:91:PHE:CE1	1:A:296:LYS:HG2	2.54	0.41
1:A:228:PHE:HB2	1:B:205:ILE:HG21	2.01	0.41
1:B:96:TYR:HE2	1:B:104:VAL:HG21	1.84	0.41
1:A:170:PRO:HB2	1:A:203:PHE:HE2	1.85	0.41
1:A:273:PHE:O	1:A:274:TYR:C	2.63	0.41
1:B:107:PRO:C	1:B:109:GLY:N	2.77	0.41
1:B:122:GLU:HB3	1:B:164:ALA:HA	2.02	0.41
1:A:211:HIS:CA	1:A:215:PRO:HG2	2.48	0.41
1:B:27:PRO:HA	1:B:102:TYR:HB3	2.02	0.41
1:B:144:SER:OG	1:B:146:PHE:CE2	2.73	0.41
1:B:179:ILE:HG13	1:B:180:PRO:HD2	2.00	0.41
1:A:57:LEU:HD23	1:A:57:LEU:O	2.21	0.41
1:A:95:LEU:C	1:A:95:LEU:CD1	2.85	0.41
1:A:104:VAL:CG2	1:A:105:PHE:N	2.83	0.41
1:A:224:GLY:C	1:A:226:LEU:N	2.79	0.41
1:A:243:THR:C	1:A:245:LYS:H	2.29	0.41
1:A:317:MET:HE2	1:A:318:VAL:CG2	2.50	0.41
1:B:95:LEU:HG	1:B:96:TYR:N	2.35	0.41
1:A:68:LEU:HD23	1:A:312:GLN:HE21	1.84	0.41
1:B:95:LEU:HD12	1:B:95:LEU:O	2.20	0.41
1:B:109:GLY:O	1:B:111:ASN:N	2.54	0.41
1:B:210:VAL:O	1:B:215:PRO:HD3	2.21	0.41
1:A:24:PHE:CD1	1:A:107:PRO:HG3	2.56	0.41
1:A:170:PRO:HB3	1:A:175:TRP:HE3	1.85	0.41
1:B:129:VAL:CG1	1:B:218:VAL:HG11	2.51	0.41
1:A:79:LEU:HD23	1:A:79:LEU:HA	1.92	0.41
1:A:269:ALA:O	1:A:270:GLY:C	2.63	0.41
1:B:166:ALA:O	1:B:206:TYR:OH	2.33	0.41
1:A:237:GLN:O	1:A:238:GLN:C	2.64	0.41
1:B:53:PRO:HG2	1:B:54:ILE:N	2.36	0.41
1:B:62:THR:HB	1:B:77:LEU:HD13	2.02	0.41
1:B:65:HIS:CE1	1:B:316:CYS:HB3	2.56	0.41
1:B:104:VAL:CG2	1:B:105:PHE:N	2.84	0.41
1:B:135:ARG:CG	1:B:254:VAL:HG21	2.50	0.41
1:B:179:ILE:CG1	1:B:180:PRO:HD2	2.50	0.41
1:B:206:TYR:OH	1:B:211:HIS:NE2	2.54	0.41
1:A:189:ILE:HB	1:A:191:TYR:HE1	1.82	0.41
1:B:48:ILE:HD12	1:B:95:LEU:CD2	2.51	0.41
1:B:53:PRO:HG2	1:B:54:ILE:H	1.86	0.41
1:A:109:GLY:O	1:A:111:ASN:N	2.54	0.40
1:A:134:GLU:HG2	1:A:148:PHE:CE1	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:224:GLY:C	1:B:226:LEU:N	2.78	0.40
1:B:305:ILE:HG23	1:B:309:MET:SD	2.61	0.40
1:A:9:PHE:CG	1:A:23:PRO:HD2	2.56	0.40
1:A:40:LEU:HD22	1:A:293:PHE:CE2	2.56	0.40
1:B:228:PHE:CE2	1:B:233:ALA:CB	2.84	0.40
1:A:48:ILE:HD12	1:A:95:LEU:HD23	2.04	0.40
1:A:122:GLU:HG3	1:A:167:CYS:HB3	2.02	0.40
1:A:220:PHE:O	1:A:223:TYR:HB3	2.22	0.40
1:A:37:PHE:CE1	1:A:185:CYS:SG	3.14	0.40
1:A:81:VAL:CG1	1:A:85:PHE:CE2	3.02	0.40
1:A:4:THR:O	1:A:9:PHE:HE1	2.05	0.40
1:A:180:PRO:HB3	1:A:184:GLN:HA	2.03	0.40
1:B:46:LEU:HA	1:B:46:LEU:HD23	1.87	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	326/349 (93%)	241 (74%)	61 (19%)	24 (7%)	1	5
1	B	326/349 (93%)	240 (74%)	63 (19%)	23 (7%)	1	6
All	All	652/698 (93%)	481 (74%)	124 (19%)	47 (7%)	1	5

All (47) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	144	SER
1	A	176	SER
1	A	177	ARG
1	A	195	HIS
1	A	230	VAL

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Mol	Chain	Res	Type
1	A	238	GLN
1	A	326	ASN
1	B	176	SER
1	B	177	ARG
1	B	195	HIS
1	B	238	GLN
1	A	142	PRO
1	A	147	ARG
1	A	182	GLY
1	A	225	GLN
1	A	323	CYS
1	B	142	PRO
1	B	150	GLU
1	B	182	GLY
1	B	200	ASN
1	B	213	ILE
1	B	230	VAL
1	B	233	ALA
1	B	323	CYS
1	B	326	ASN
1	A	108	THR
1	A	110	CYS
1	A	190	ASP
1	A	200	ASN
1	A	213	ILE
1	B	105	PHE
1	B	110	CYS
1	B	144	SER
1	B	225	GLN
1	B	231	LYS
1	A	105	PHE
1	A	231	LYS
1	A	233	ALA
1	B	108	THR
1	B	53	PRO
1	B	147	ARG
1	A	53	PRO
1	A	162	VAL
1	A	300	VAL
1	B	300	VAL
1	A	227	VAL
1	B	227	VAL

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	278/296 (94%)	262 (94%)	16 (6%)	18	45
1	B	278/296 (94%)	258 (93%)	20 (7%)	13	39
All	All	556/592 (94%)	520 (94%)	36 (6%)	15	42

All (36) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	39	MET
1	A	50	LEU
1	A	95	LEU
1	A	97	THR
1	A	108	THR
1	A	110	CYS
1	A	113	GLU
1	A	125	LEU
1	A	172	LEU
1	A	193	THR
1	A	206	TYR
1	A	244	GLN
1	A	255	ILE
1	A	263	ILE
1	A	266	LEU
1	A	275	ILE
1	B	39	MET
1	B	50	LEU
1	B	95	LEU
1	B	97	THR
1	B	108	THR
1	B	110	CYS
1	B	113	GLU
1	B	125	LEU
1	B	140	CYS
1	B	172	LEU
1	B	193	THR

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Mol	Chain	Res	Type
1	B	201	GLU
1	B	206	TYR
1	B	244	GLN
1	B	255	ILE
1	B	263	ILE
1	B	266	LEU
1	B	275	ILE
1	B	323	CYS
1	B	325	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	ASN
1	A	200	ASN
1	A	312	GLN
1	B	73	ASN
1	B	312	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

4 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	NAG	B	1327	1	14,14,15	2.59	5 (35%)	17,19,21	1.42	4 (23%)
3	NAG	A	1327	1	14,14,15	2.48	5 (35%)	17,19,21	1.46	3 (17%)
2	RET	B	1296	1	20,20,21	1.44	2 (10%)	27,27,28	1.73	7 (25%)
2	RET	A	1296	1	20,20,21	1.49	3 (15%)	27,27,28	1.70	5 (18%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NAG	B	1327	1	-	3/6/23/26	0/1/1/1
3	NAG	A	1327	1	-	3/6/23/26	0/1/1/1
2	RET	B	1296	1	-	1/13/30/31	0/1/1/1
2	RET	A	1296	1	-	1/13/30/31	0/1/1/1

All (15) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	B	1327	NAG	C1-C2	6.88	1.61	1.52
3	A	1327	NAG	C1-C2	5.75	1.60	1.52
3	A	1327	NAG	C3-C2	4.29	1.61	1.52
2	A	1296	RET	C14-C13	4.11	1.36	1.33
3	B	1327	NAG	O5-C1	4.07	1.50	1.43
2	B	1296	RET	C14-C13	3.96	1.36	1.33
3	A	1327	NAG	O5-C1	3.54	1.49	1.43
3	B	1327	NAG	C3-C2	3.32	1.59	1.52
2	B	1296	RET	C1-C6	3.08	1.57	1.53
2	A	1296	RET	C1-C6	3.00	1.57	1.53
3	A	1327	NAG	O5-C5	2.70	1.48	1.43
3	B	1327	NAG	O5-C5	2.68	1.48	1.43
3	B	1327	NAG	C2-N2	2.46	1.50	1.46
3	A	1327	NAG	C8-C7	2.31	1.55	1.50
2	A	1296	RET	C4-C5	2.02	1.54	1.51

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1296	RET	C20-C13-C12	4.18	124.48	118.09
2	A	1296	RET	C8-C9-C10	-3.72	113.16	119.01
2	B	1296	RET	C20-C13-C12	3.66	123.68	118.09
2	B	1296	RET	C8-C9-C10	-3.52	113.48	119.01
2	B	1296	RET	C7-C8-C9	3.01	130.68	126.23
2	A	1296	RET	C11-C12-C13	2.83	134.12	126.36
2	B	1296	RET	C19-C9-C10	2.80	127.35	122.82
2	B	1296	RET	C11-C12-C13	2.79	134.01	126.36
3	B	1327	NAG	O5-C5-C6	2.74	113.00	107.66
3	A	1327	NAG	C1-C2-N2	-2.68	106.21	110.43
2	A	1296	RET	C19-C9-C10	2.61	127.05	122.82
2	A	1296	RET	C7-C8-C9	2.52	129.96	126.23
3	B	1327	NAG	O5-C1-C2	2.42	115.03	111.29
2	B	1296	RET	C18-C5-C6	2.34	127.03	124.48
2	B	1296	RET	C10-C11-C12	2.31	129.89	123.20
3	B	1327	NAG	O3-C3-C2	2.28	114.13	109.40
3	A	1327	NAG	O3-C3-C2	2.09	113.73	109.40
3	B	1327	NAG	C3-C4-C5	-2.06	106.49	110.23
3	A	1327	NAG	O5-C5-C6	2.06	111.66	107.66

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	B	1327	NAG	O5-C5-C6-O6
3	B	1327	NAG	C4-C5-C6-O6
3	A	1327	NAG	C4-C5-C6-O6
3	A	1327	NAG	O5-C5-C6-O6
3	A	1327	NAG	C3-C2-N2-C7
3	B	1327	NAG	C3-C2-N2-C7
2	A	1296	RET	C10-C11-C12-C13
2	B	1296	RET	C10-C11-C12-C13

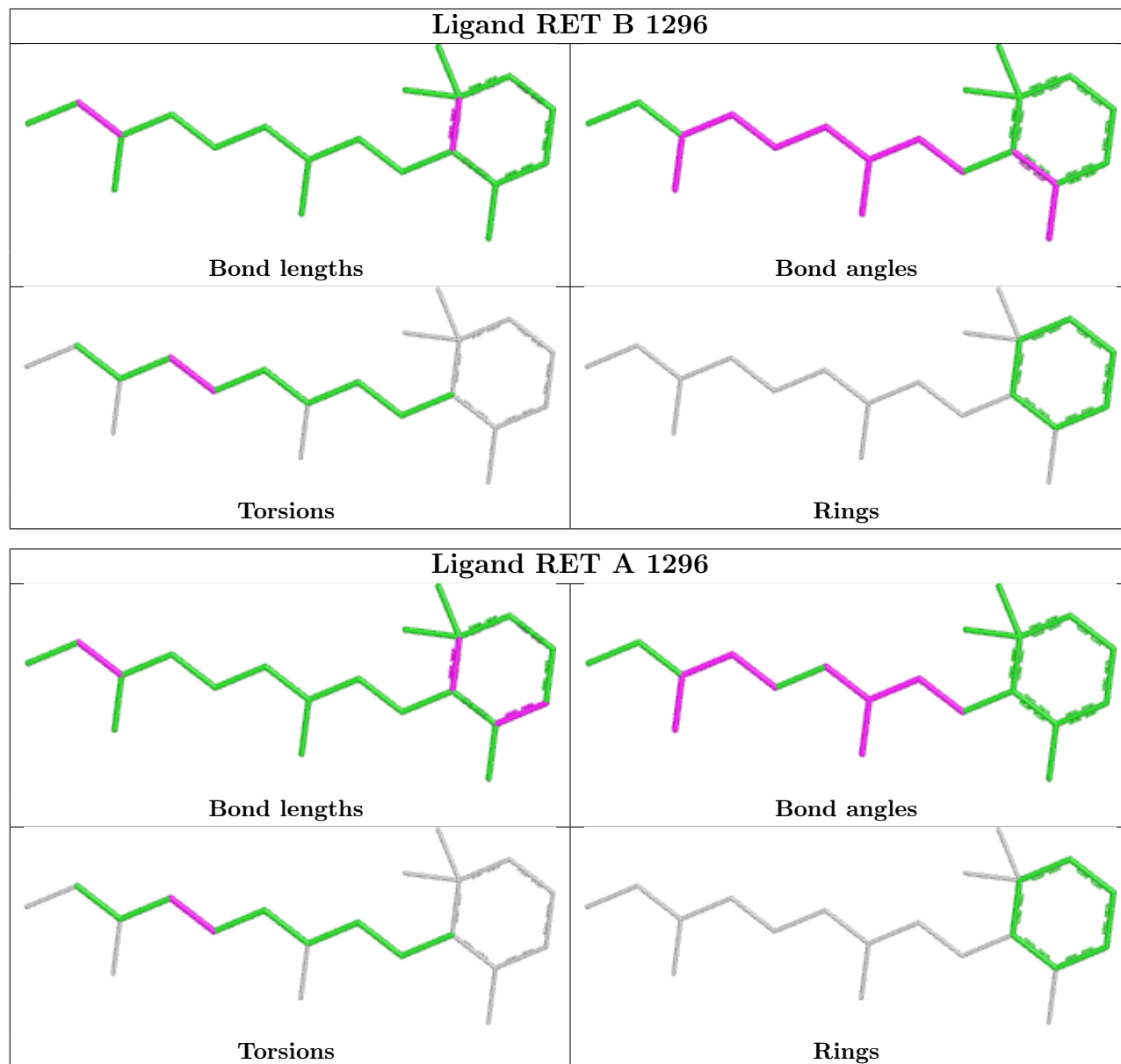
There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	1327	NAG	1	0
2	B	1296	RET	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

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



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues

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	327/349 (93%)	-1.22	0 100 100	10, 44, 112, 160	0
1	B	327/349 (93%)	-1.19	0 100 100	12, 42, 115, 154	0
All	All	654/698 (93%)	-1.20	0 100 100	10, 43, 115, 160	0

There are no RSRZ outliers to report.

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

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

6.3 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

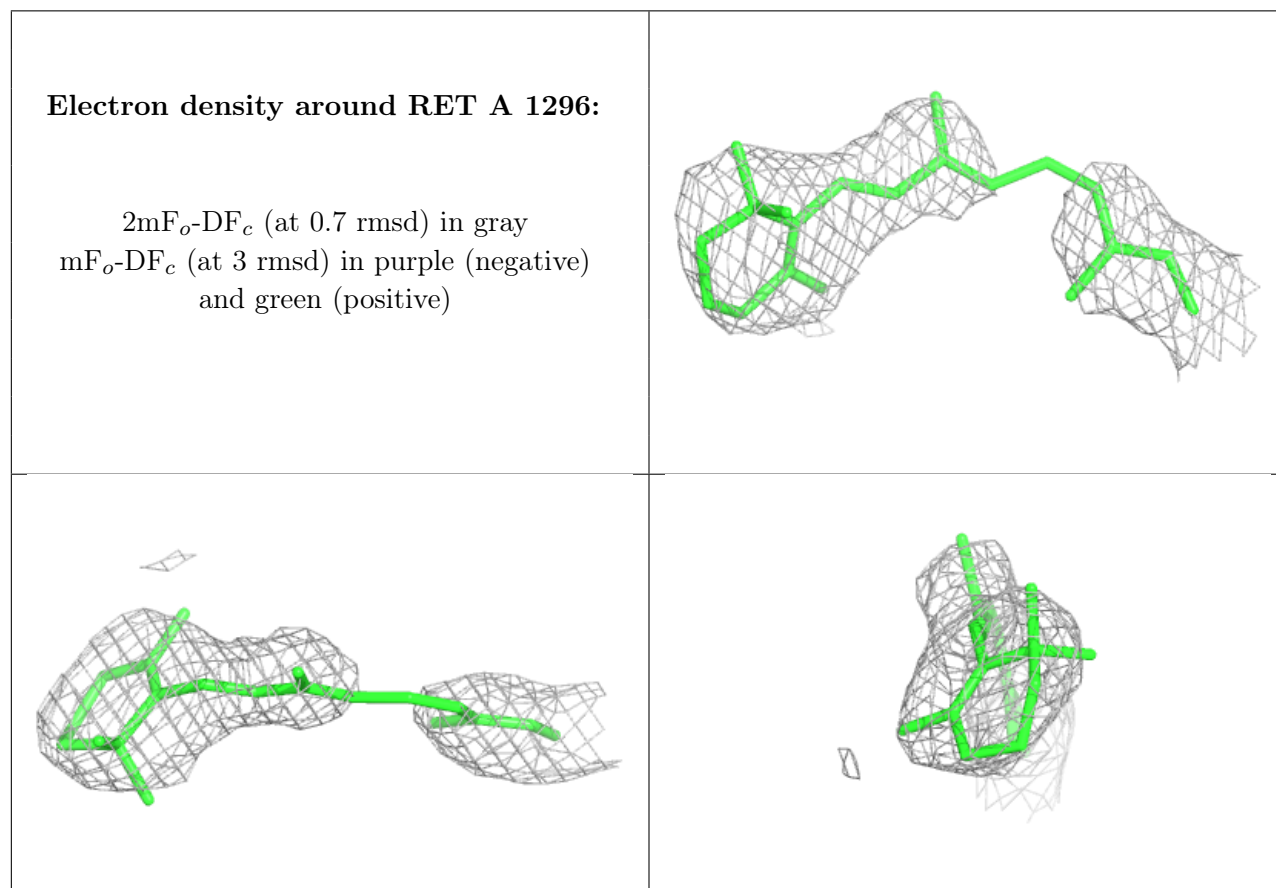
6.4 Ligands [i](#)

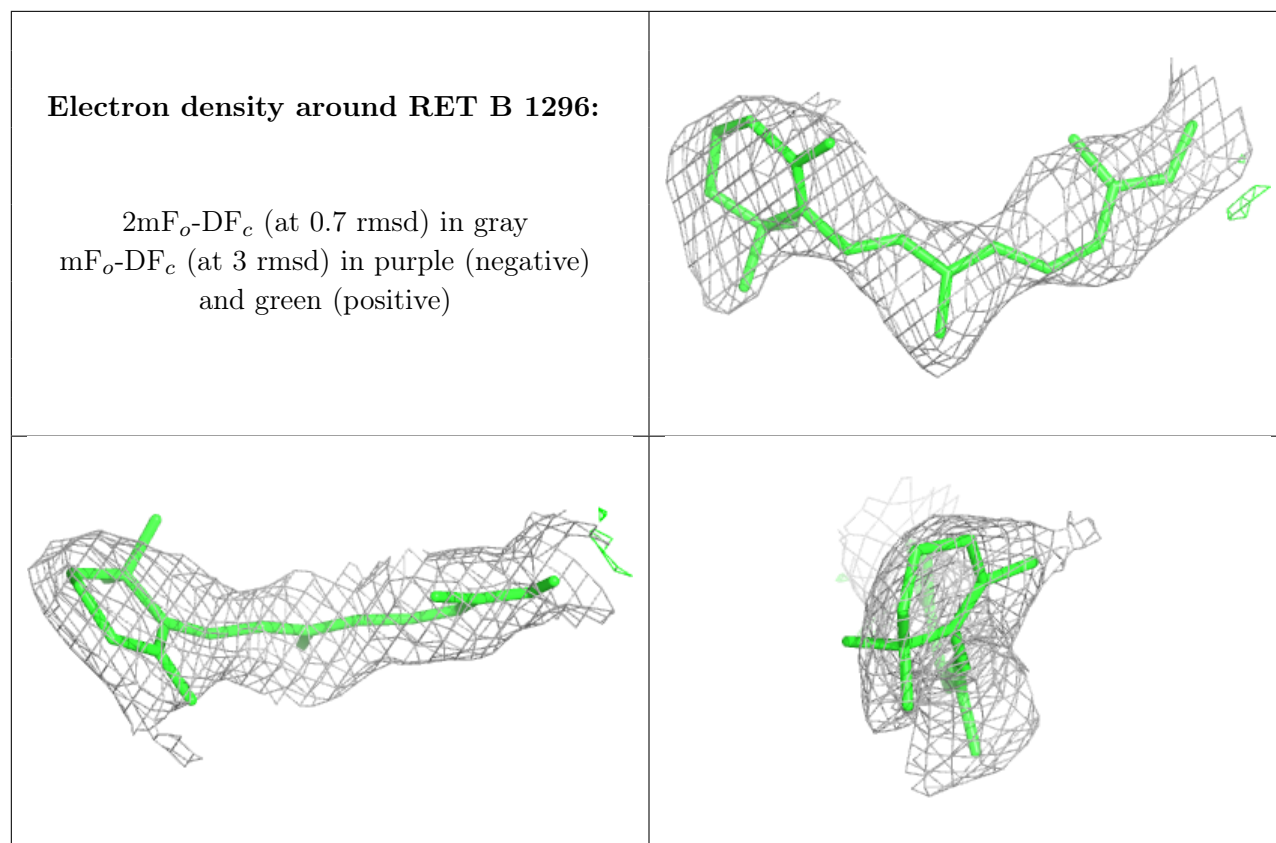
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q< 0.9’ lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	NAG	A	1327	14/15	0.97	0.05	58,63,64,65	0
3	NAG	B	1327	14/15	0.98	0.05	50,53,54,54	0
2	RET	A	1296	20/21	0.99	0.06	35,38,41,41	0
2	RET	B	1296	20/21	0.99	0.04	34,37,41,41	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers

as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





6.5 Other polymers [i](#)

There are no such residues in this entry.