



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 11:57 AM UTC

PDB ID : 1MQS / pdb_00001mq
Title : Crystal structure of Sly1p in complex with an N-terminal peptide of Sed5p
Authors : Bracher, A.; Weissenhorn, W.
Deposited on : 2002-09-17
Resolution : 3.00 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

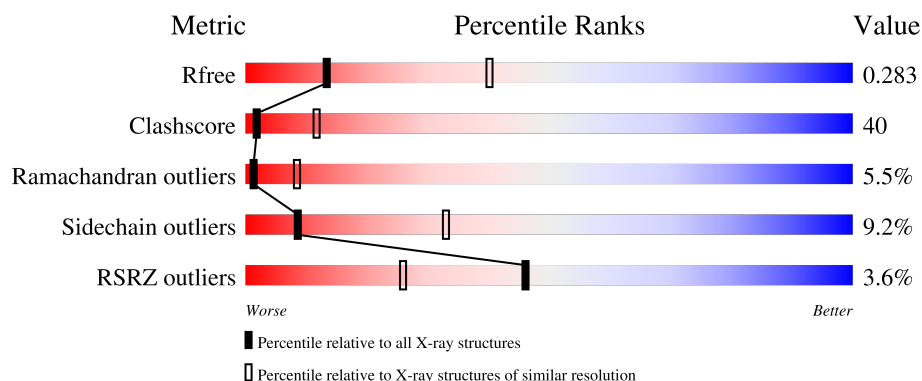
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

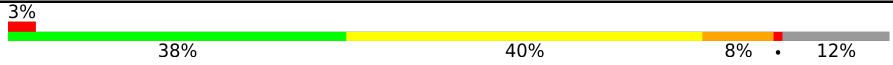
The reported resolution of this entry is 3.00 Å.

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	2672 (3.00-3.00)
Clashscore	190562	2977 (3.00-3.00)
Ramachandran outliers	187476	2877 (3.00-3.00)
Sidechain outliers	187428	2880 (3.00-3.00)
RSRZ outliers	180081	2671 (3.00-3.00)

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	671	
2	B	50	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4725 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 Sly1 Protein.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	588	Total	C	N	O	S	Se	0	0	0
			4506	2871	731	892	2	10			

There are 15 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-4	GLY	-	cloning artifact	UNP P22213
A	-3	LYS	-	cloning artifact	UNP P22213
A	-2	SER	-	cloning artifact	UNP P22213
A	-1	ALA	-	cloning artifact	UNP P22213
A	0	SER	-	cloning artifact	UNP P22213
A	17	MSE	MET	modified residue	UNP P22213
A	25	MSE	MET	modified residue	UNP P22213
A	273	MSE	MET	modified residue	UNP P22213
A	283	MSE	MET	modified residue	UNP P22213
A	329	MSE	MET	modified residue	UNP P22213
A	401	MSE	MET	modified residue	UNP P22213
A	497	MSE	MET	modified residue	UNP P22213
A	502	MSE	MET	modified residue	UNP P22213
A	566	MSE	MET	modified residue	UNP P22213
A	637	MSE	MET	modified residue	UNP P22213

- Molecule 2 is a protein called Integral Membrane Protein SED5.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	26	Total	C	N	O	S	0	0	0
			199	122	38	38	1			

There are 5 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	-4	GLY	-	cloning artifact	UNP Q01590

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-3	ALA	-	cloning artifact	UNP Q01590
B	-2	MET	-	cloning artifact	UNP Q01590
B	-1	ALA	-	cloning artifact	UNP Q01590
B	0	GLY	-	cloning artifact	UNP Q01590

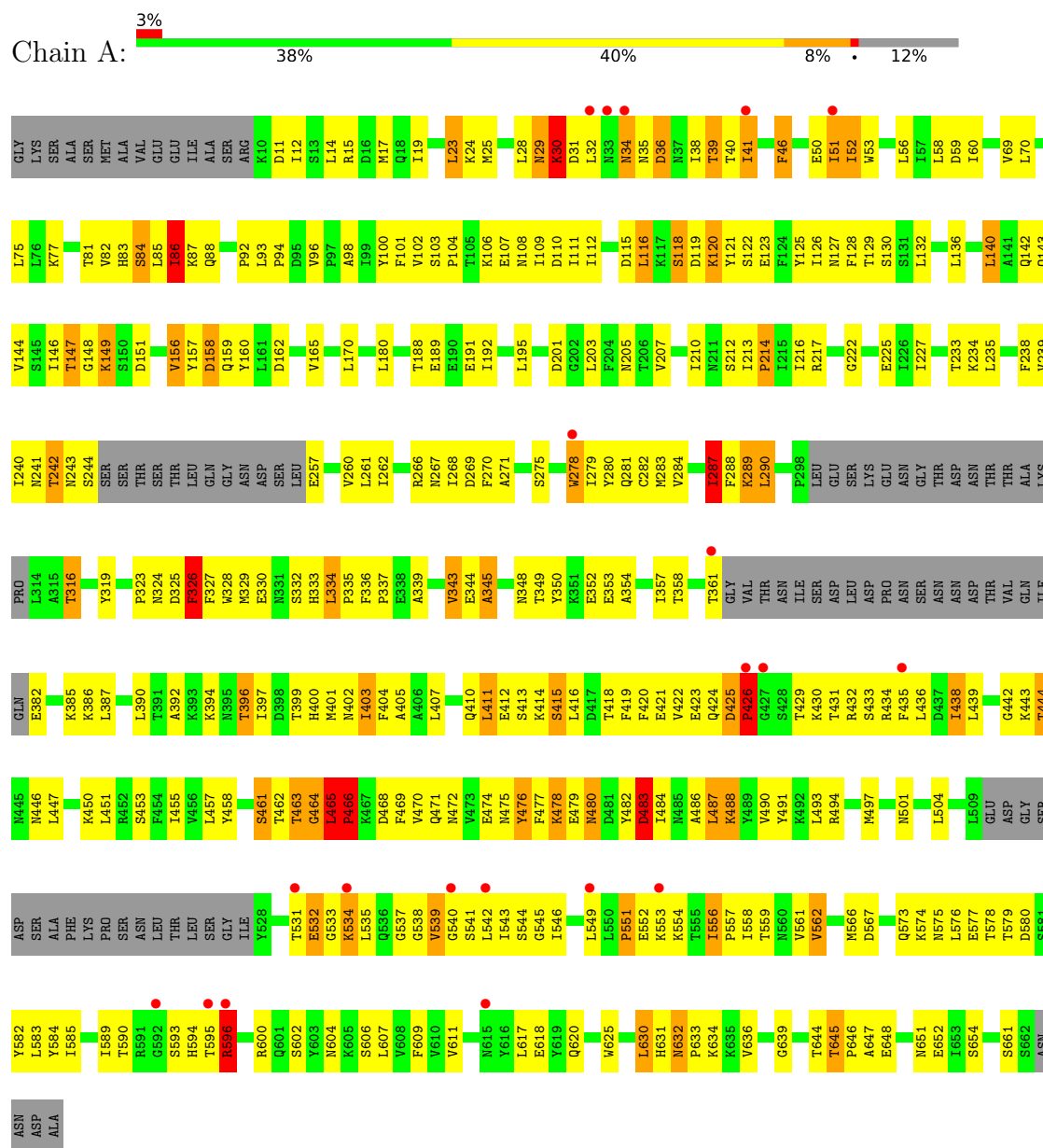
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	19	Total O 19 19	0	0
3	B	1	Total O 1 1	0	0

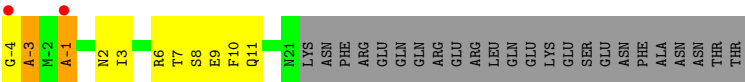
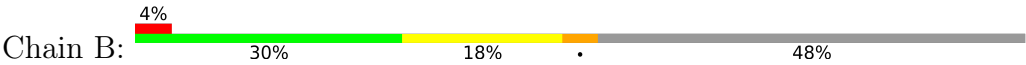
3 Residue-property plots

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: Sly1 Protein



• Molecule 2: Integral Membrane Protein SED5



4 Data and refinement statistics

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants a, b, c, α , β , γ	161.06Å 161.06Å 88.21Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 3.00 30.00 – 3.00	Depositor EDS
% Data completeness (in resolution range)	99.9 (30.00-3.00) 99.8 (30.00-3.00)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.38 (at 2.90Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.257 , 0.290 0.255 , 0.283	Depositor DCC
R_{free} test set	1328 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	83.8	Xtriage
Anisotropy	0.767	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 71.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.52$, $\langle L^2 \rangle = 0.36$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4725	wwPDB-VP
Average B, all atoms (Å ²)	99.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.02% 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

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	0.58	2/4574 (0.0%)	1.14	35/6213 (0.6%)
2	B	0.56	0/200	1.04	1/264 (0.4%)
All	All	0.58	2/4774 (0.0%)	1.14	36/6477 (0.6%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	30	LYS	CA-CB	-7.69	1.40	1.53
1	A	30	LYS	C-O	-7.54	1.14	1.24

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	596	ARG	N-CA-C	-11.67	93.55	109.54
1	A	403	ILE	N-CA-C	-10.78	99.84	110.72
1	A	38	ILE	N-CA-C	-10.67	95.80	108.82
1	A	594	HIS	N-CA-C	-10.32	99.00	111.69
1	A	532	GLU	N-CA-C	-8.48	100.89	111.75
1	A	189	GLU	N-CA-C	-7.80	103.49	113.16
1	A	396	THR	N-CA-C	-6.89	103.70	111.07
1	A	480	ASN	N-CA-C	-6.81	104.96	113.20
1	A	112	ILE	N-CA-C	-6.75	103.90	110.72
1	A	567	ASP	CA-C-N	6.57	125.80	118.97
1	A	567	ASP	C-N-CA	6.57	125.80	118.97
1	A	158	ASP	N-CA-C	-6.34	98.88	108.96
1	A	465	LEU	CA-C-N	6.33	127.75	119.84
1	A	465	LEU	C-N-CA	6.33	127.75	119.84
1	A	34	ASN	N-CA-C	6.14	116.35	108.24
1	A	630	LEU	N-CA-C	-6.11	100.61	110.32
1	A	602	SER	N-CA-C	-5.97	101.17	110.42
1	A	537	GLY	N-CA-C	-5.94	107.06	115.43
1	A	87	LYS	N-CA-C	5.90	119.74	112.54
1	A	149	LYS	N-CA-C	5.88	120.06	112.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	287	ILE	N-CA-C	5.84	118.39	111.09
1	A	222	GLY	CA-C-N	5.78	127.06	119.84
1	A	222	GLY	C-N-CA	5.78	127.06	119.84
1	A	316	THR	N-CA-C	5.77	117.53	107.49
2	B	-1	ALA	N-CA-C	-5.75	106.30	113.38
1	A	29	ASN	N-CA-C	-5.70	104.38	111.90
1	A	147	THR	N-CA-C	-5.66	101.06	110.17
1	A	12	ILE	N-CA-C	5.64	116.33	108.89
1	A	148	GLY	N-CA-C	-5.30	106.49	115.34
1	A	345	ALA	N-CA-C	-5.28	105.98	112.90
1	A	86	ILE	CB-CA-C	-5.24	105.07	112.14
1	A	51	ILE	N-CA-C	5.20	117.04	108.97
1	A	582	TYR	N-CA-C	5.17	117.17	109.25
1	A	534	LYS	N-CA-C	5.16	121.79	110.80
1	A	488	LYS	N-CA-C	-5.14	104.60	111.24
1	A	46	PHE	N-CA-C	-5.08	105.19	111.33

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	4506	0	4353	362	0
2	B	199	0	199	12	0
3	A	19	0	0	2	0
3	B	1	0	0	0	0
All	All	4725	0	4552	367	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 40.

All (367) 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:429:THR:HA	1:A:432:ARG:HB2	1.26	1.12
1:A:539:VAL:HG12	1:A:540:GLY:H	1.09	1.11
1:A:576:LEU:HD22	1:A:600:ARG:HH11	1.09	1.11
1:A:562:VAL:HG12	1:A:566:MSE:HE2	1.27	1.09
1:A:29:ASN:HB3	1:A:32:LEU:HD12	1.35	1.07
1:A:283:MSE:HE1	1:A:336:PHE:HB2	1.06	1.03
1:A:414:LYS:HE3	1:A:416:LEU:HD13	1.39	1.01
1:A:25:MSE:HE2	1:A:127:ASN:HB3	1.40	1.00
1:A:504:LEU:HD23	1:A:578:THR:HG21	1.42	1.00
1:A:24:LYS:HE2	1:A:32:LEU:HG	1.43	0.97
1:A:470:VAL:O	1:A:474:GLU:HB2	1.65	0.96
1:A:539:VAL:HG12	1:A:540:GLY:N	1.77	0.96
1:A:645:THR:HB	1:A:648:GLU:HG3	1.44	0.94
1:A:283:MSE:HE1	1:A:336:PHE:CB	1.97	0.94
1:A:539:VAL:CG1	1:A:540:GLY:H	1.81	0.94
1:A:216:ILE:HB	1:A:585:ILE:HG23	1.49	0.92
1:A:562:VAL:CG1	1:A:566:MSE:HE2	1.97	0.92
1:A:576:LEU:HD22	1:A:600:ARG:NH1	1.87	0.89
1:A:279:ILE:HD11	1:A:282:CYS:SG	2.13	0.88
1:A:539:VAL:HG13	1:A:543:ILE:HB	1.56	0.88
1:A:52:ILE:HD13	1:A:52:ILE:H	1.39	0.87
1:A:283:MSE:CE	1:A:336:PHE:HB2	2.00	0.87
1:A:283:MSE:O	1:A:287:ILE:HG12	1.74	0.87
1:A:327:PHE:CZ	1:A:343:VAL:HG23	2.10	0.86
1:A:158:ASP:OD2	2:B:6:ARG:HD2	1.76	0.86
1:A:539:VAL:HG12	1:A:541:SER:H	1.43	0.84
1:A:424:GLN:HE21	1:A:551:PRO:HG2	1.42	0.82
1:A:549:LEU:O	1:A:551:PRO:HD3	1.78	0.82
1:A:14:LEU:HA	1:A:17:MSE:HE3	1.62	0.81
1:A:266:ARG:O	1:A:266:ARG:HD3	1.81	0.81
1:A:261:LEU:HD12	1:A:607:LEU:HB3	1.63	0.80
1:A:25:MSE:HE2	1:A:127:ASN:CB	2.12	0.80
1:A:461:SER:O	1:A:462:THR:HG22	1.82	0.79
1:A:25:MSE:CE	1:A:127:ASN:HB3	2.13	0.79
1:A:235:LEU:O	1:A:239:VAL:HG23	1.83	0.78
1:A:336:PHE:HB3	1:A:337:PRO:HD3	1.64	0.78
1:A:278:TRP:O	1:A:283:MSE:HE2	1.84	0.78
1:A:336:PHE:CE1	1:A:407:LEU:HD23	2.19	0.78
1:A:418:THR:O	1:A:421:GLU:HG2	1.83	0.78
1:A:261:LEU:HD13	1:A:607:LEU:HD23	1.67	0.77
1:A:504:LEU:CD2	1:A:578:THR:HG21	2.15	0.77
1:A:556:ILE:HG23	1:A:557:PRO:HD2	1.68	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:29:ASN:HB3	1:A:32:LEU:CD1	2.14	0.74
1:A:104:PRO:HD3	1:A:129:THR:HG22	1.68	0.74
1:A:558:ILE:O	1:A:562:VAL:HG23	1.88	0.74
1:A:432:ARG:HB3	1:A:469:PHE:HZ	1.53	0.72
1:A:486:ALA:HB2	1:A:654:SER:HB3	1.72	0.72
1:A:539:VAL:HG12	1:A:541:SER:N	2.03	0.72
1:A:92:PRO:O	1:A:93:LEU:HD22	1.89	0.72
1:A:147:THR:HG22	1:A:149:LYS:H	1.54	0.72
1:A:109:ILE:HD11	1:A:132:LEU:HD11	1.71	0.71
1:A:278:TRP:HA	1:A:278:TRP:CE3	2.25	0.70
1:A:201:ASP:OD1	1:A:234:LYS:HE2	1.91	0.70
1:A:589:ILE:HG22	1:A:590:THR:N	2.06	0.70
1:A:281:GLN:HG3	1:A:332:SER:O	1.91	0.70
1:A:213:ILE:HD13	1:A:239:VAL:HG22	1.71	0.70
1:A:278:TRP:HA	1:A:278:TRP:HE3	1.56	0.69
1:A:30:LYS:C	1:A:32:LEU:H	2.01	0.69
1:A:93:LEU:HD12	1:A:96:VAL:HG21	1.75	0.68
1:A:425:ASP:N	1:A:426:PRO:HD2	2.09	0.68
1:A:419:PHE:O	1:A:423:GLU:HG3	1.95	0.67
1:A:52:ILE:H	1:A:52:ILE:CD1	2.07	0.67
1:A:386:LYS:O	1:A:387:LEU:HD23	1.94	0.67
1:A:432:ARG:HD3	1:A:469:PHE:CE1	2.29	0.67
1:A:350:TYR:CE2	1:A:394:LYS:HG3	2.29	0.67
1:A:645:THR:CB	1:A:648:GLU:HG3	2.21	0.66
1:A:425:ASP:HB3	1:A:431:THR:OG1	1.97	0.65
1:A:645:THR:HB	1:A:648:GLU:CG	2.22	0.65
1:A:436:LEU:HD12	1:A:436:LEU:H	1.60	0.65
1:A:596:ARG:NH1	1:A:596:ARG:HG2	2.10	0.65
1:A:339:ALA:HB1	1:A:404:PHE:HZ	1.62	0.65
1:A:422:VAL:HG12	1:A:457:LEU:HD11	1.78	0.65
1:A:35:ASN:O	1:A:36:ASP:HB2	1.95	0.65
1:A:350:TYR:CZ	1:A:394:LYS:HG3	2.32	0.65
1:A:589:ILE:HG22	1:A:590:THR:H	1.60	0.65
1:A:84:SER:HB3	1:A:88:GLN:OE1	1.98	0.64
1:A:596:ARG:CG	1:A:596:ARG:HH11	2.09	0.64
1:A:288:PHE:CE1	1:A:407:LEU:HD12	2.32	0.64
1:A:458:TYR:CE2	1:A:491:TYR:HA	2.33	0.64
1:A:429:THR:HA	1:A:432:ARG:CB	2.16	0.64
1:A:484:ILE:HG22	1:A:487:LEU:HB2	1.80	0.63
1:A:59:ASP:HB2	1:A:86:ILE:HD11	1.79	0.63
1:A:69:VAL:HG23	1:A:70:LEU:HD13	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:477:PHE:HB3	1:A:484:ILE:HD11	1.81	0.63
1:A:279:ILE:HD13	1:A:620:GLN:CG	2.28	0.63
1:A:270:PHE:CD2	1:A:493:LEU:HD13	2.34	0.63
1:A:436:LEU:HA	1:A:439:LEU:HG	1.79	0.63
1:A:596:ARG:HG2	1:A:596:ARG:HH11	1.64	0.63
1:A:556:ILE:HG23	1:A:557:PRO:CD	2.29	0.62
1:A:447:LEU:HG	1:A:482:TYR:CD2	2.35	0.62
1:A:573:GLN:O	1:A:577:GLU:HG3	1.99	0.62
1:A:59:ASP:CB	1:A:86:ILE:HD11	2.29	0.62
1:A:188:THR:HB	1:A:191:GLU:H	1.64	0.62
1:A:103:SER:HA	1:A:129:THR:HG22	1.82	0.62
1:A:288:PHE:O	1:A:289:LYS:C	2.43	0.62
1:A:132:LEU:O	2:B:6:ARG:NH1	2.33	0.62
1:A:279:ILE:HD13	1:A:620:GLN:HG2	1.80	0.62
1:A:53:TRP:HB3	1:A:81:THR:OG1	2.01	0.61
1:A:539:VAL:HG13	1:A:543:ILE:CB	2.28	0.61
1:A:561:VAL:HG13	1:A:579:THR:HG22	1.83	0.61
1:A:217:ARG:HD2	1:A:584:TYR:HA	1.83	0.61
1:A:432:ARG:HB3	1:A:469:PHE:CZ	2.33	0.61
1:A:420:PHE:O	1:A:424:GLN:HG2	2.01	0.61
1:A:279:ILE:HD13	1:A:620:GLN:CD	2.27	0.60
1:A:280:TYR:O	1:A:284:VAL:HG23	2.02	0.60
1:A:81:THR:HG22	1:A:82:VAL:HG23	1.84	0.60
1:A:336:PHE:HE1	1:A:407:LEU:HD23	1.67	0.60
1:A:325:ASP:O	1:A:327:PHE:N	2.35	0.59
1:A:539:VAL:HG11	1:A:544:SER:H	1.66	0.59
1:A:576:LEU:CD2	1:A:600:ARG:HH11	2.01	0.59
1:A:556:ILE:HG22	1:A:558:ILE:HG22	1.84	0.59
1:A:34:ASN:CG	1:A:35:ASN:H	2.09	0.59
1:A:631:HIS:O	1:A:632:ASN:HB2	2.02	0.59
1:A:267:ASN:HA	1:A:270:PHE:CE1	2.38	0.59
1:A:425:ASP:O	1:A:426:PRO:C	2.46	0.58
1:A:25:MSE:HE3	1:A:159:GLN:CD	2.27	0.58
1:A:531:THR:C	1:A:533:GLY:N	2.57	0.58
1:A:424:GLN:NE2	1:A:551:PRO:HG2	2.16	0.58
1:A:574:LYS:O	1:A:578:THR:HG22	2.04	0.58
1:A:576:LEU:HB3	1:A:600:ARG:NH1	2.18	0.58
1:A:140:LEU:HD13	2:B:10:PHE:CE1	2.38	0.58
1:A:348:ASN:O	1:A:352:GLU:HG2	2.04	0.58
1:A:278:TRP:C	1:A:283:MSE:HE2	2.28	0.58
2:B:7:THR:O	2:B:11:GLN:HG3	2.03	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:129:THR:HG23	1:A:130:SER:N	2.19	0.57
1:A:227:ILE:HD12	1:A:611:VAL:HG21	1.85	0.57
1:A:458:TYR:HE2	1:A:491:TYR:HA	1.69	0.57
1:A:25:MSE:HE3	1:A:159:GLN:OE1	2.04	0.57
1:A:397:ILE:O	1:A:401:MSE:HB2	2.04	0.57
1:A:484:ILE:HG22	1:A:484:ILE:O	2.04	0.57
1:A:210:ILE:CD1	1:A:607:LEU:HD22	2.34	0.57
1:A:238:PHE:HE2	2:B:-4:GLY:HA3	1.70	0.57
1:A:566:MSE:HE1	1:A:636:VAL:CG2	2.35	0.57
1:A:644:THR:HG23	1:A:648:GLU:HB2	1.87	0.56
1:A:243:ASN:O	1:A:244:SER:HB2	2.04	0.56
1:A:630:LEU:O	1:A:631:HIS:HB2	2.06	0.56
1:A:281:GLN:HE21	1:A:620:GLN:NE2	2.03	0.56
1:A:358:THR:HA	3:A:673:HOH:O	2.04	0.56
1:A:106:LYS:HE3	1:A:110:ASP:OD2	2.05	0.56
1:A:103:SER:HA	1:A:129:THR:CG2	2.37	0.55
1:A:494:ARG:HH11	1:A:494:ARG:HG3	1.71	0.55
1:A:19:ILE:HG12	1:A:69:VAL:HB	1.87	0.55
1:A:279:ILE:CD1	1:A:282:CYS:SG	2.93	0.55
1:A:56:LEU:O	1:A:83:HIS:HA	2.07	0.55
1:A:271:ALA:HB2	1:A:497:MSE:HE2	1.88	0.55
1:A:323:PRO:O	1:A:324:ASN:HB2	2.06	0.55
1:A:93:LEU:N	1:A:94:PRO:HD3	2.23	0.54
1:A:25:MSE:HE1	1:A:101:PHE:CD1	2.42	0.54
1:A:401:MSE:HE3	1:A:401:MSE:HA	1.89	0.54
1:A:425:ASP:N	1:A:426:PRO:CD	2.70	0.54
1:A:160:TYR:HA	1:A:205:ASN:ND2	2.22	0.54
1:A:645:THR:HG22	1:A:647:ALA:N	2.23	0.54
1:A:147:THR:HG22	1:A:149:LYS:N	2.22	0.54
1:A:461:SER:OG	1:A:494:ARG:NH1	2.41	0.54
1:A:25:MSE:HE2	1:A:127:ASN:CG	2.32	0.54
1:A:461:SER:O	1:A:462:THR:CG2	2.56	0.53
1:A:25:MSE:HE3	1:A:159:GLN:CG	2.39	0.53
1:A:69:VAL:O	1:A:70:LEU:HD12	2.07	0.53
1:A:100:TYR:O	1:A:127:ASN:HB2	2.08	0.53
1:A:19:ILE:HG23	1:A:70:LEU:HD11	1.91	0.53
1:A:216:ILE:HB	1:A:585:ILE:CG2	2.29	0.53
1:A:279:ILE:CD1	1:A:620:GLN:HG2	2.38	0.53
1:A:539:VAL:HG13	1:A:543:ILE:H	1.72	0.53
1:A:28:LEU:HD13	1:A:125:TYR:CZ	2.44	0.53
1:A:392:ALA:O	1:A:396:THR:HG23	2.09	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:645:THR:HG22	1:A:647:ALA:H	1.73	0.53
1:A:207:VAL:HA	1:A:210:ILE:HD12	1.91	0.52
1:A:358:THR:O	1:A:358:THR:HG22	2.09	0.52
1:A:482:TYR:O	1:A:483:ASP:C	2.52	0.52
1:A:539:VAL:HG21	1:A:544:SER:OG	2.09	0.52
1:A:403:ILE:O	1:A:407:LEU:HB2	2.08	0.52
1:A:354:ALA:CB	1:A:390:LEU:HD21	2.40	0.52
1:A:419:PHE:HE2	1:A:450:LYS:HA	1.74	0.52
1:A:316:THR:HG23	1:A:316:THR:O	2.08	0.52
1:A:343:VAL:HG12	1:A:344:GLU:N	2.25	0.52
1:A:539:VAL:CG1	1:A:540:GLY:N	2.48	0.52
1:A:69:VAL:C	1:A:70:LEU:HD12	2.35	0.52
1:A:463:THR:HG23	1:A:463:THR:O	2.09	0.52
1:A:39:THR:O	1:A:39:THR:CG2	2.57	0.52
1:A:128:PHE:CE1	1:A:132:LEU:HD22	2.46	0.51
1:A:533:GLY:O	1:A:535:LEU:N	2.33	0.51
1:A:418:THR:OG1	1:A:419:PHE:N	2.44	0.51
1:A:539:VAL:HG11	1:A:544:SER:OG	2.10	0.51
1:A:435:PHE:O	1:A:438:ILE:HG22	2.10	0.51
1:A:576:LEU:HB3	1:A:600:ARG:HH12	1.75	0.51
1:A:633:PRO:O	1:A:634:LYS:HD3	2.10	0.51
1:A:24:LYS:HE2	1:A:32:LEU:CG	2.30	0.51
1:A:444:THR:C	1:A:446:ASN:H	2.19	0.51
1:A:589:ILE:CG2	1:A:590:THR:N	2.75	0.51
1:A:344:GLU:OE2	1:A:532:GLU:O	2.29	0.50
1:A:238:PHE:CE2	2:B:-4:GLY:HA3	2.45	0.50
1:A:288:PHE:O	1:A:290:LEU:N	2.45	0.50
1:A:29:ASN:C	1:A:30:LYS:O	2.54	0.50
1:A:102:VAL:HG12	1:A:103:SER:N	2.27	0.50
1:A:464:GLY:O	1:A:465:LEU:CB	2.60	0.50
1:A:58:LEU:O	1:A:85:LEU:HA	2.12	0.50
1:A:283:MSE:HB3	1:A:287:ILE:HD11	1.93	0.50
1:A:484:ILE:HG21	1:A:487:LEU:HD12	1.93	0.50
1:A:339:ALA:HB1	1:A:404:PHE:CZ	2.44	0.50
1:A:327:PHE:CD2	1:A:400:HIS:CD2	3.01	0.49
1:A:240:ILE:C	1:A:242:THR:H	2.18	0.49
1:A:538:GLY:C	1:A:539:VAL:HG23	2.38	0.49
1:A:589:ILE:CG2	1:A:590:THR:H	2.26	0.49
1:A:29:ASN:O	1:A:32:LEU:HB2	2.13	0.49
1:A:30:LYS:C	1:A:32:LEU:N	2.69	0.49
1:A:109:ILE:CD1	1:A:136:LEU:HB3	2.43	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:162:ASP:H	1:A:205:ASN:ND2	2.11	0.49
1:A:432:ARG:HE	1:A:466:PRO:HD2	1.78	0.49
1:A:439:LEU:C	1:A:439:LEU:HD12	2.38	0.49
1:A:645:THR:HG22	1:A:647:ALA:HB3	1.95	0.49
1:A:575:ASN:HA	1:A:578:THR:CG2	2.43	0.48
1:A:86:ILE:HG12	1:A:108:ASN:ND2	2.28	0.48
1:A:270:PHE:CZ	1:A:493:LEU:HD22	2.49	0.48
1:A:429:THR:O	1:A:430:LYS:C	2.55	0.48
1:A:477:PHE:CB	1:A:484:ILE:HD11	2.42	0.48
1:A:102:VAL:HG11	1:A:108:ASN:CG	2.38	0.48
1:A:325:ASP:HB3	1:A:328:TRP:HB3	1.95	0.48
1:A:562:VAL:HG12	1:A:566:MSE:CE	2.19	0.48
1:A:328:TRP:C	1:A:330:GLU:H	2.20	0.48
1:A:432:ARG:HH21	1:A:466:PRO:HD3	1.77	0.48
1:A:472:ASN:HA	1:A:475:ASN:HD22	1.77	0.48
1:A:281:GLN:HE21	1:A:620:GLN:HE21	1.62	0.48
1:A:225:GLU:HG3	1:A:583:LEU:HD22	1.96	0.48
1:A:262:ILE:HD12	1:A:562:VAL:HG13	1.95	0.47
1:A:39:THR:O	1:A:39:THR:HG23	2.13	0.47
1:A:281:GLN:NE2	1:A:620:GLN:HE21	2.12	0.47
1:A:414:LYS:O	1:A:415:SER:C	2.57	0.47
1:A:30:LYS:O	1:A:32:LEU:N	2.47	0.47
1:A:386:LYS:C	1:A:387:LEU:HD23	2.40	0.47
1:A:486:ALA:C	1:A:488:LYS:H	2.21	0.47
1:A:539:VAL:HG13	1:A:543:ILE:N	2.29	0.47
1:A:261:LEU:CD1	1:A:607:LEU:HB3	2.37	0.47
1:A:455:ILE:HD13	1:A:490:VAL:HG21	1.97	0.47
1:A:240:ILE:O	1:A:242:THR:N	2.38	0.47
1:A:86:ILE:N	1:A:86:ILE:HD12	2.30	0.47
1:A:290:LEU:C	1:A:290:LEU:HD22	2.40	0.47
1:A:319:TYR:CD2	1:A:399:THR:HG23	2.48	0.47
1:A:575:ASN:HA	1:A:578:THR:HG22	1.96	0.47
2:B:-1:ALA:O	2:B:2:ASN:HB2	2.14	0.47
1:A:162:ASP:H	1:A:205:ASN:HD22	1.62	0.47
1:A:278:TRP:NE1	1:A:546:ILE:HG23	2.30	0.47
1:A:213:ILE:CD1	1:A:239:VAL:HG22	2.41	0.47
1:A:242:THR:HG22	1:A:243:ASN:ND2	2.29	0.47
1:A:609:PHE:HA	1:A:639:GLY:O	2.15	0.47
1:A:631:HIS:O	1:A:632:ASN:CB	2.63	0.47
1:A:644:THR:CG2	1:A:648:GLU:HB2	2.45	0.47
1:A:129:THR:CG2	1:A:130:SER:N	2.77	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:396:THR:O	1:A:399:THR:HB	2.15	0.47
1:A:102:VAL:CG1	1:A:103:SER:N	2.77	0.46
1:A:104:PRO:HD3	1:A:129:THR:CG2	2.41	0.46
1:A:115:ASP:OD1	1:A:120:LYS:HE2	2.15	0.46
1:A:462:THR:O	1:A:463:THR:HB	2.14	0.46
1:A:645:THR:HG23	1:A:646:PRO:HD2	1.96	0.46
1:A:645:THR:O	1:A:646:PRO:C	2.58	0.46
1:A:102:VAL:CG1	1:A:108:ASN:CG	2.89	0.46
1:A:349:THR:O	1:A:353:GLU:HB2	2.15	0.46
1:A:484:ILE:CG2	1:A:487:LEU:HD12	2.45	0.46
1:A:140:LEU:HD22	1:A:144:VAL:HG23	1.98	0.46
1:A:645:THR:CG2	1:A:647:ALA:HB3	2.46	0.46
1:A:410:GLN:O	1:A:414:LYS:HG2	2.15	0.46
1:A:107:GLU:O	1:A:111:ILE:HG13	2.16	0.45
1:A:217:ARG:HD2	1:A:217:ARG:HA	1.79	0.45
1:A:354:ALA:HB1	1:A:390:LEU:HD21	1.96	0.45
1:A:645:THR:CG2	1:A:647:ALA:H	2.29	0.45
1:A:156:VAL:HG13	2:B:7:THR:HA	1.99	0.45
1:A:126:ILE:O	1:A:156:VAL:HA	2.16	0.45
1:A:212:SER:O	1:A:214:PRO:HD3	2.16	0.45
1:A:382:GLU:N	3:A:684:HOH:O	2.49	0.45
1:A:160:TYR:HA	1:A:205:ASN:HD21	1.81	0.45
1:A:541:SER:O	1:A:542:LEU:HD23	2.17	0.45
1:A:539:VAL:CG1	1:A:541:SER:H	2.22	0.45
1:A:411:LEU:C	1:A:411:LEU:HD12	2.41	0.45
1:A:217:ARG:NH1	1:A:217:ARG:HG3	2.32	0.45
1:A:336:PHE:CD1	1:A:407:LEU:HD23	2.51	0.45
1:A:486:ALA:O	1:A:488:LYS:N	2.49	0.45
1:A:281:GLN:NE2	1:A:620:GLN:NE2	2.65	0.45
1:A:419:PHE:HB3	1:A:453:SER:OG	2.17	0.45
1:A:486:ALA:C	1:A:488:LYS:N	2.75	0.45
1:A:325:ASP:O	1:A:326:PHE:C	2.60	0.44
1:A:651:ASN:O	1:A:652:GLU:C	2.60	0.44
1:A:41:ILE:H	1:A:41:ILE:HG13	1.37	0.44
1:A:217:ARG:HG3	1:A:217:ARG:HH11	1.82	0.44
1:A:217:ARG:CD	1:A:584:TYR:HA	2.46	0.44
1:A:104:PRO:HG2	1:A:130:SER:OG	2.17	0.44
1:A:201:ASP:OD1	1:A:234:LYS:CE	2.61	0.44
1:A:466:PRO:C	1:A:468:ASP:N	2.75	0.44
1:A:118:SER:O	1:A:119:ASP:HB2	2.17	0.44
1:A:419:PHE:HD2	1:A:453:SER:OG	2.01	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:566:MSE:HE1	1:A:636:VAL:HG22	1.98	0.44
1:A:477:PHE:CG	1:A:484:ILE:HD11	2.52	0.44
1:A:451:LEU:O	1:A:455:ILE:HG13	2.17	0.44
1:A:29:ASN:O	1:A:32:LEU:HD12	2.18	0.44
1:A:270:PHE:O	1:A:271:ALA:C	2.61	0.44
1:A:345:ALA:O	1:A:349:THR:HG23	2.18	0.44
1:A:19:ILE:CG1	1:A:69:VAL:HB	2.48	0.44
1:A:116:LEU:HD22	1:A:116:LEU:O	2.18	0.44
1:A:122:SER:O	1:A:123:GLU:HG3	2.17	0.44
1:A:15:ARG:HB2	1:A:165:VAL:HG11	1.99	0.43
1:A:32:LEU:HD13	1:A:157:TYR:OH	2.18	0.43
1:A:268:ILE:HG12	1:A:268:ILE:O	2.17	0.43
1:A:558:ILE:HG23	1:A:559:THR:N	2.33	0.43
1:A:584:TYR:C	1:A:584:TYR:CD2	2.95	0.43
1:A:25:MSE:HE3	1:A:159:GLN:HG2	1.99	0.43
1:A:59:ASP:C	1:A:59:ASP:OD1	2.61	0.43
1:A:160:TYR:C	1:A:205:ASN:HD22	2.26	0.43
1:A:556:ILE:HG23	1:A:557:PRO:N	2.33	0.43
1:A:438:ILE:HD13	1:A:438:ILE:O	2.19	0.43
1:A:539:VAL:HG11	1:A:544:SER:N	2.32	0.43
1:A:34:ASN:CG	1:A:35:ASN:N	2.76	0.43
1:A:343:VAL:HG11	1:A:401:MSE:CE	2.49	0.43
1:A:143:GLN:O	1:A:146:ILE:HG12	2.18	0.43
1:A:422:VAL:HG11	1:A:435:PHE:HD1	1.83	0.43
1:A:46:PHE:CE2	1:A:51:ILE:HB	2.54	0.43
1:A:354:ALA:O	1:A:358:THR:OG1	2.20	0.43
1:A:257:GLU:N	1:A:604:ASN:HD21	2.17	0.43
1:A:471:GLN:O	1:A:475:ASN:ND2	2.52	0.43
1:A:279:ILE:HD13	1:A:620:GLN:NE2	2.34	0.42
1:A:402:ASN:HA	1:A:405:ALA:HB3	2.01	0.42
1:A:552:GLU:O	1:A:554:LYS:N	2.50	0.42
1:A:158:ASP:OD2	2:B:6:ARG:CD	2.59	0.42
1:A:447:LEU:HG	1:A:482:TYR:CE2	2.54	0.42
2:B:7:THR:O	2:B:10:PHE:HB3	2.19	0.42
1:A:35:ASN:O	1:A:36:ASP:CB	2.65	0.42
1:A:344:GLU:O	1:A:348:ASN:HB2	2.19	0.42
1:A:385:LYS:HD3	1:A:386:LYS:H	1.84	0.42
1:A:59:ASP:O	1:A:60:ILE:C	2.62	0.42
1:A:266:ARG:HH22	1:A:618:GLU:CD	2.27	0.42
1:A:15:ARG:CB	1:A:165:VAL:HG11	2.50	0.42
2:B:8:SER:O	2:B:9:GLU:C	2.61	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:279:ILE:HA	1:A:335:PRO:HA	2.01	0.42
1:A:334:LEU:HA	1:A:335:PRO:HD3	1.88	0.42
2:B:-3:ALA:C	2:B:-1:ALA:H	2.28	0.42
1:A:288:PHE:CE1	1:A:407:LEU:CD1	3.01	0.42
1:A:327:PHE:HB3	1:A:400:HIS:CD2	2.54	0.42
1:A:333:HIS:CE1	1:A:334:LEU:HD13	2.54	0.42
1:A:439:LEU:HB2	1:A:476:TYR:CE2	2.54	0.42
1:A:484:ILE:CG2	1:A:487:LEU:HB2	2.48	0.42
1:A:436:LEU:HA	1:A:439:LEU:CG	2.46	0.42
1:A:412:GLU:C	1:A:414:LYS:H	2.28	0.41
1:A:425:ASP:H	1:A:426:PRO:CD	2.32	0.41
1:A:463:THR:O	1:A:465:LEU:N	2.53	0.41
1:A:188:THR:OG1	1:A:191:GLU:HG3	2.20	0.41
1:A:556:ILE:CG2	1:A:558:ILE:HG22	2.49	0.41
1:A:584:TYR:O	1:A:596:ARG:HB3	2.20	0.41
1:A:23:LEU:HA	1:A:23:LEU:HD12	1.70	0.41
1:A:416:LEU:C	1:A:418:THR:H	2.27	0.41
1:A:461:SER:O	1:A:463:THR:HG22	2.20	0.41
1:A:566:MSE:HE3	1:A:625:TRP:CH2	2.55	0.41
1:A:566:MSE:SE	1:A:606:SER:HB3	2.70	0.41
1:A:260:VAL:O	1:A:606:SER:HA	2.21	0.41
1:A:478:LYS:H	1:A:478:LYS:HG2	1.69	0.41
1:A:545:GLY:O	1:A:549:LEU:HG	2.21	0.41
1:A:270:PHE:CE2	1:A:493:LEU:HD13	2.56	0.41
1:A:432:ARG:CB	1:A:469:PHE:CZ	3.03	0.41
1:A:484:ILE:C	1:A:486:ALA:N	2.79	0.41
1:A:497:MSE:O	1:A:501:ASN:HB2	2.21	0.41
1:A:539:VAL:CG1	1:A:544:SER:H	2.32	0.41
1:A:479:GLU:HG3	1:A:480:ASN:N	2.36	0.40
1:A:538:GLY:C	1:A:539:VAL:CG2	2.94	0.40
1:A:85:LEU:HB2	1:A:88:GLN:HG3	2.03	0.40
1:A:203:LEU:HD23	1:A:203:LEU:HA	1.89	0.40
1:A:323:PRO:O	1:A:324:ASN:CB	2.69	0.40
1:A:433:SER:C	1:A:435:PHE:N	2.79	0.40
1:A:357:ILE:HG21	1:A:387:LEU:HD13	2.02	0.40
1:A:539:VAL:CG1	1:A:543:ILE:H	2.35	0.40
1:A:466:PRO:C	1:A:468:ASP:H	2.28	0.40
1:A:98:ALA:HB2	1:A:121:TYR:CG	2.57	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	578/671 (86%)	467 (81%)	79 (14%)	32 (6%)	1	8
2	B	24/50 (48%)	20 (83%)	3 (12%)	1 (4%)	2	13
All	All	602/721 (84%)	487 (81%)	82 (14%)	33 (6%)	1	8

All (33) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	30	LYS
1	A	31	ASP
1	A	36	ASP
1	A	275	SER
1	A	326	PHE
1	A	426	PRO
1	A	442	GLY
1	A	461	SER
1	A	464	GLY
1	A	534	LYS
1	A	40	THR
1	A	289	LYS
2	B	-3	ALA
1	A	50	GLU
1	A	241	ASN
1	A	415	SER
1	A	434	ARG
1	A	444	THR
1	A	463	THR
1	A	483	ASP
1	A	487	LEU
1	A	593	SER
1	A	118	SER
1	A	269	ASP
1	A	413	SER

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Mol	Chain	Res	Type
1	A	443	LYS
1	A	465	LEU
1	A	425	ASP
1	A	551	PRO
1	A	466	PRO
1	A	476	TYR
1	A	539	VAL
1	A	632	ASN

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	491/597 (82%)	445 (91%)	46 (9%)	8	32
2	B	20/45 (44%)	19 (95%)	1 (5%)	22	56
All	All	511/642 (80%)	464 (91%)	47 (9%)	8	33

All (47) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	11	ASP
1	A	23	LEU
1	A	39	THR
1	A	41	ILE
1	A	52	ILE
1	A	75	LEU
1	A	77	LYS
1	A	84	SER
1	A	86	ILE
1	A	116	LEU
1	A	120	LYS
1	A	140	LEU
1	A	142	GLN
1	A	151	ASP
1	A	156	VAL

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Mol	Chain	Res	Type
1	A	170	LEU
1	A	180	LEU
1	A	192	ILE
1	A	195	LEU
1	A	214	PRO
1	A	233	THR
1	A	242	THR
1	A	278	TRP
1	A	287	ILE
1	A	290	LEU
1	A	326	PHE
1	A	329	MSE
1	A	334	LEU
1	A	343	VAL
1	A	361	THR
1	A	411	LEU
1	A	426	PRO
1	A	438	ILE
1	A	465	LEU
1	A	466	PRO
1	A	478	LYS
1	A	483	ASP
1	A	553	LYS
1	A	556	ILE
1	A	562	VAL
1	A	580	ASP
1	A	595	THR
1	A	596	ARG
1	A	617	LEU
1	A	645	THR
1	A	661	SER
2	B	3	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	37	ASN
1	A	108	ASN
1	A	135	ASN
1	A	142	GLN
1	A	205	ASN
1	A	243	ASN

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Mol	Chain	Res	Type
1	A	276	HIS
1	A	281	GLN
1	A	331	ASN
1	A	342	ASN
1	A	424	GLN
1	A	445	ASN
1	A	620	GLN
1	A	632	ASN
1	A	651	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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](#)

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	578/671 (86%)	0.17	20 (3%)	47 27	57, 94, 159, 177	0
2	B	26/50 (52%)	0.34	2 (7%)	19 10	67, 81, 138, 144	0
All	All	604/721 (83%)	0.18	22 (3%)	46 26	57, 94, 159, 177	0

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	278	TRP	3.7
1	A	51	ILE	3.7
1	A	361	THR	3.7
1	A	32	LEU	3.6
1	A	33	ASN	3.3
1	A	549	LEU	3.1
1	A	41	ILE	3.1
1	A	595	THR	2.9
1	A	553	LYS	2.9
2	B	-4	GLY	2.9
1	A	534	LYS	2.9
1	A	34	ASN	2.8
1	A	542	LEU	2.7
1	A	615	ASN	2.6
2	B	-1	ALA	2.5
1	A	596	ARG	2.4
1	A	531	THR	2.4
1	A	435	PHE	2.2
1	A	540	GLY	2.1
1	A	426	PRO	2.1
1	A	592	GLY	2.1
1	A	427	GLY	2.0

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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.