



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 10, 2026 – 12:08 AM UTC

PDB ID : 1YO6 / pdb_00001yo6
Title : Crystal Structure of the putative Carbonyl Reductase Sniffer of *Caenorhabditis elegans*
Authors : Southeast Collaboratory for Structural Genomics (SECSG)
Deposited on : 2005-01-26
Resolution : 2.60 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
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

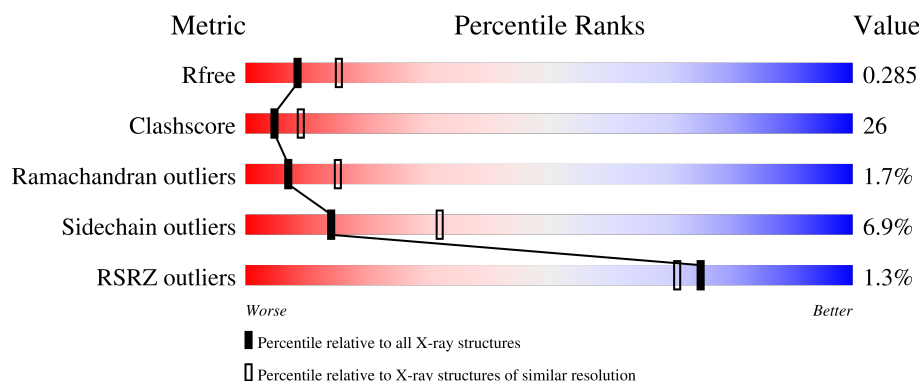
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	250	57% 29% 7% 5%
1	B	250	58% 30% 6% 6%
1	C	250	56% 32% 7% 5%
1	D	250	50% 38% 7% 6%
1	E	250	50% 39% 5% 6%

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Mol	Chain	Length	Quality of chain
1	F	250	 A horizontal bar chart showing the quality of chain F. The bar is divided into four segments: a small red segment at the beginning labeled '4%', a large green segment labeled '48%', a large yellow segment labeled '40%', and a small orange segment at the end labeled '8%'. The segments are separated by thin black lines.

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 10878 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 Putative Carbonyl Reductase Sniffer.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	237	Total	C	N	O	S	0	0	0
			1786	1121	311	348	6			
1	B	235	Total	C	N	O	S	0	0	0
			1765	1105	308	346	6			
1	C	240	Total	C	N	O	S	0	0	0
			1809	1135	315	353	6			
1	D	236	Total	C	N	O	S	0	0	0
			1779	1116	310	347	6			
1	E	235	Total	C	N	O	S	0	0	0
			1764	1105	308	346	5			
1	F	230	Total	C	N	O	S	0	0	0
			1739	1091	303	340	5			

- Molecule 2 is water.

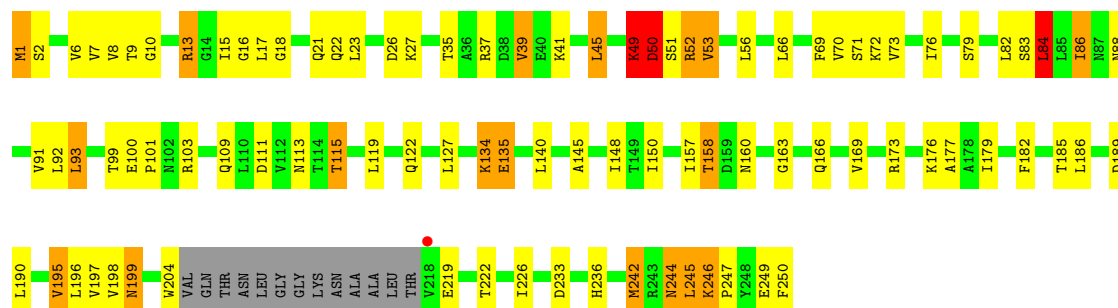
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	51	Total	O	0	0
			51	51		
2	B	40	Total	O	0	0
			40	40		
2	C	50	Total	O	0	0
			50	50		
2	D	35	Total	O	0	0
			35	35		
2	E	31	Total	O	0	0
			31	31		
2	F	29	Total	O	0	0
			29	29		

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.

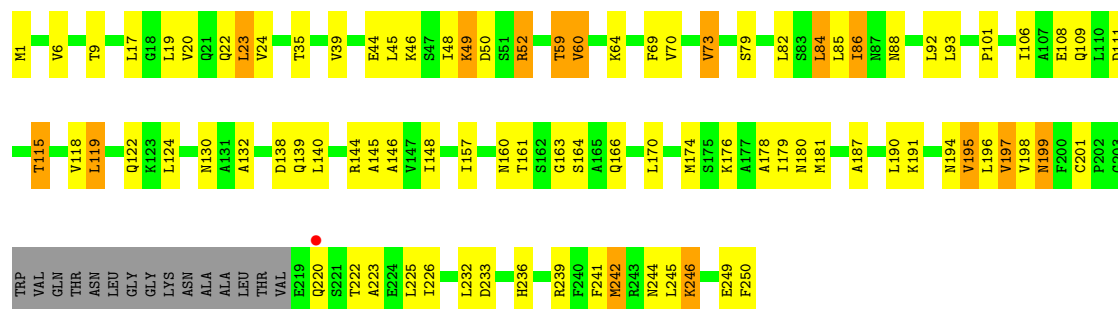
• Molecule 1: Putative Carbonyl Reductase Sniffer

Chain A: 



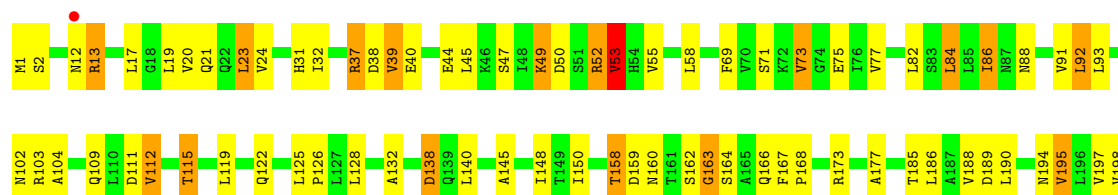
• Molecule 1: Putative Carbonyl Reductase Sniffer

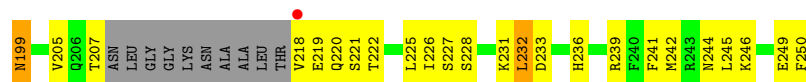
Chain B: 



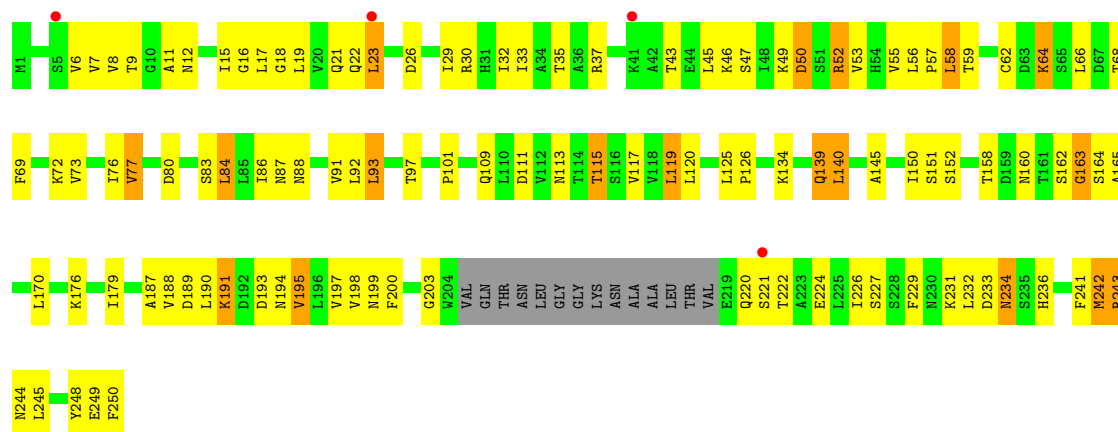
• Molecule 1: Putative Carbonyl Reductase Sniffer

Chain C: 

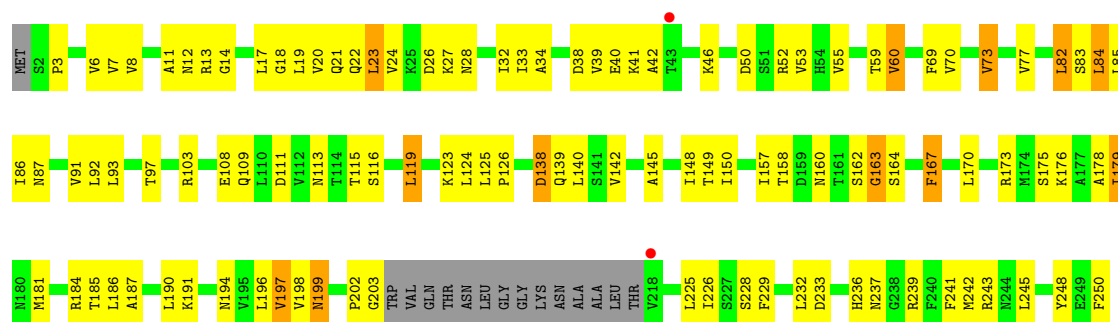




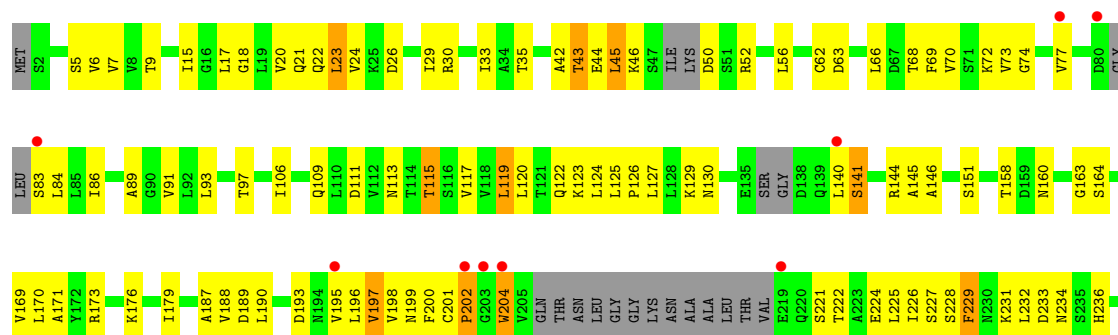
• Molecule 1: Putative Carbonyl Reductase Sniffer



• Molecule 1: Putative Carbonyl Reductase Sniffer



• Molecule 1: Putative Carbonyl Reductase Sniffer



R239	F240	F241	M242	R243	N244	L245		Y248	E249	F250
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4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	74.87Å 86.32Å 242.56Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	47.12 – 2.60 47.12 – 2.60	Depositor EDS
% Data completeness (in resolution range)	93.4 (47.12-2.60) 93.4 (47.12-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	7.56 (at 2.61Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.218 , 0.286 0.218 , 0.285	Depositor DCC
R_{free} test set	2307 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	48.0	Xtriage
Anisotropy	0.048	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 55.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10878	wwPDB-VP
Average B, all atoms (Å ²)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.20% 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 [i](#)

5.1 Standard geometry [i](#)

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	0.49	0/1807	0.99	10/2446 (0.4%)
1	B	0.50	0/1784	0.93	5/2413 (0.2%)
1	C	0.50	0/1830	0.99	9/2478 (0.4%)
1	D	0.46	0/1800	0.97	5/2436 (0.2%)
1	E	0.43	0/1783	0.94	4/2413 (0.2%)
1	F	0.43	0/1757	0.96	6/2376 (0.3%)
All	All	0.47	0/10761	0.96	39/14562 (0.3%)

There are no bond length outliers.

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	F	202	PRO	N-CA-C	-9.40	97.28	111.57
1	A	219	GLU	N-CA-C	-8.05	103.44	113.18
1	D	242	MET	N-CA-C	-8.01	98.00	110.42
1	C	242	MET	N-CA-C	-7.29	100.02	110.59
1	A	177	ALA	N-CA-C	-6.67	103.93	111.14
1	A	51	SER	N-CA-C	6.65	120.49	112.38
1	B	246	LYS	N-CA-C	-6.27	101.75	109.65
1	F	229	PHE	N-CA-C	-6.25	104.56	111.82
1	D	245	LEU	N-CA-C	6.06	120.00	112.24
1	C	177	ALA	N-CA-C	-6.00	104.82	111.36
1	B	201	CYS	N-CA-C	-5.89	99.32	109.15
1	A	247	PRO	N-CA-C	5.78	120.41	111.21
1	B	180	ASN	N-CA-C	-5.75	105.14	111.82
1	A	244	ASN	N-CA-C	-5.62	106.31	112.72
1	F	201	CYS	CA-C-N	-5.57	114.85	120.31
1	F	201	CYS	C-N-CA	-5.57	114.85	120.31
1	A	84	LEU	N-CA-C	5.55	117.46	108.41
1	F	242	MET	N-CA-C	-5.51	102.37	110.52
1	B	60	VAL	N-CA-C	5.49	120.77	109.34
1	C	219	GLU	N-CA-C	-5.47	106.11	112.89
1	D	77	VAL	N-CA-C	5.43	116.92	111.00

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	113	ASN	N-CA-C	5.42	119.16	112.54
1	C	112	VAL	N-CA-C	5.39	116.17	110.72
1	A	169	VAL	N-CA-C	5.31	117.51	111.88
1	F	164	SER	N-CA-C	-5.30	105.09	112.30
1	A	45	LEU	N-CA-C	-5.26	105.45	111.14
1	A	242	MET	N-CA-C	-5.26	102.74	110.52
1	A	37	ARG	N-CA-C	-5.26	105.46	111.14
1	D	203	GLY	N-CA-C	-5.18	104.02	112.83
1	E	167	PHE	CA-C-N	5.14	126.26	119.84
1	E	167	PHE	C-N-CA	5.14	126.26	119.84
1	C	246	LYS	CA-C-N	5.13	125.05	119.76
1	C	246	LYS	C-N-CA	5.13	125.05	119.76
1	C	232	LEU	N-CA-C	5.12	117.74	110.10
1	E	60	VAL	N-CA-C	5.10	118.20	111.17
1	B	242	MET	N-CA-C	-5.08	103.22	110.59
1	C	52	ARG	N-CA-C	-5.08	106.93	113.23
1	E	33	ILE	N-CA-C	5.07	114.58	106.88
1	C	53	VAL	N-CA-C	5.00	115.29	107.99

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	1786	0	1840	87	0
1	B	1765	0	1821	80	0
1	C	1809	0	1864	97	0
1	D	1779	0	1831	117	0
1	E	1764	0	1818	93	0
1	F	1739	0	1779	109	0
2	A	51	0	0	3	0
2	B	40	0	0	4	0
2	C	50	0	0	3	0
2	D	35	0	0	1	0
2	E	31	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	F	29	0	0	0	0
All	All	10878	0	10953	553	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 26.

All (553) 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:158:THR:HG21	1:C:249:GLU:HB3	1.45	0.93
1:E:111:ASP:HA	1:E:115:THR:HG22	1.52	0.92
1:F:231:LYS:HE2	1:F:231:LYS:HA	1.52	0.91
1:A:50:ASP:OD1	1:A:52:ARG:HG2	1.71	0.91
1:C:163:GLY:HA2	1:C:173:ARG:HE	1.37	0.90
1:B:93:LEU:H	1:B:109:GLN:NE2	1.71	0.89
1:E:111:ASP:HA	1:E:115:THR:CG2	2.03	0.88
1:D:15:ILE:HD12	1:D:15:ILE:H	1.38	0.87
1:B:160:ASN:HD21	1:B:163:GLY:H	1.23	0.87
1:B:233:ASP:H	1:B:236:HIS:HD2	1.20	0.85
1:A:111:ASP:HA	1:A:115:THR:HG23	1.59	0.85
1:D:158:THR:HG23	1:D:250:PHE:O	1.78	0.84
1:F:127:LEU:HD12	1:F:127:LEU:H	1.41	0.84
1:C:158:THR:HG22	1:C:250:PHE:OXT	1.77	0.84
1:C:233:ASP:H	1:C:236:HIS:HD2	1.22	0.83
1:B:111:ASP:HA	1:B:115:THR:HG23	1.60	0.83
1:B:93:LEU:H	1:B:109:GLN:HE22	1.22	0.82
1:E:19:LEU:O	1:E:23:LEU:HD22	1.78	0.81
1:E:163:GLY:HA2	1:E:173:ARG:HE	1.44	0.81
1:E:138:ASP:HA	1:E:194:ASN:ND2	1.96	0.81
1:C:93:LEU:H	1:C:109:GLN:HE22	1.27	0.81
1:A:1:MET:HA	1:A:1:MET:HE2	1.63	0.81
1:A:15:ILE:HD12	1:A:15:ILE:H	1.45	0.81
1:B:1:MET:HE3	1:B:1:MET:HA	1.63	0.80
1:E:138:ASP:HA	1:E:194:ASN:HD21	1.47	0.80
1:A:166:GLN:HG3	1:C:205:VAL:HG23	1.63	0.80
1:D:50:ASP:OD1	1:D:52:ARG:HG2	1.83	0.78
1:B:111:ASP:HA	1:B:115:THR:CG2	2.14	0.77
1:E:163:GLY:HA2	1:E:173:ARG:NE	1.99	0.77
1:B:17:LEU:HD21	1:B:44:GLU:OE2	1.85	0.77
1:D:6:VAL:HG13	1:D:84:LEU:HD13	1.66	0.77
1:E:158:THR:HG22	1:F:250:PHE:HD2	1.50	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:50:ASP:OD1	1:B:52:ARG:HG2	1.84	0.76
1:C:218:VAL:N	1:C:221:SER:HG	1.84	0.76
1:C:239:ARG:HD2	2:C:276:HOH:O	1.86	0.76
1:D:232:LEU:HA	1:D:236:HIS:CD2	2.21	0.75
1:C:93:LEU:H	1:C:109:GLN:NE2	1.85	0.74
1:F:6:VAL:HG13	1:F:84:LEU:HD22	1.68	0.74
1:D:26:ASP:OD2	1:D:29:ILE:HG13	1.86	0.74
1:D:233:ASP:H	1:D:236:HIS:HD2	1.35	0.74
1:A:158:THR:HG22	1:A:250:PHE:OXT	1.88	0.74
1:A:160:ASN:HD21	1:A:163:GLY:H	1.33	0.73
1:D:190:LEU:HB3	1:D:195:VAL:HG13	1.71	0.73
1:D:111:ASP:HA	1:D:115:THR:HG23	1.70	0.72
1:F:233:ASP:H	1:F:236:HIS:HD2	1.37	0.72
1:A:6:VAL:HG13	1:A:84:LEU:HB3	1.71	0.72
1:E:12:ASN:ND2	1:E:13:ARG:HG2	2.04	0.72
1:A:245:LEU:O	1:A:246:LYS:HB3	1.90	0.71
1:D:64:LYS:HE3	1:D:64:LYS:HA	1.71	0.71
1:F:127:LEU:HD12	1:F:127:LEU:N	2.05	0.71
1:D:115:THR:HG21	2:D:271:HOH:O	1.89	0.71
1:F:187:ALA:HA	1:F:197:VAL:HG11	1.71	0.71
1:C:158:THR:CG2	1:C:249:GLU:HB3	2.19	0.71
1:C:13:ARG:NH2	1:C:88:ASN:HD21	1.89	0.71
1:C:50:ASP:HB3	1:C:53:VAL:CG1	2.21	0.70
1:F:111:ASP:HA	1:F:115:THR:HG23	1.74	0.70
1:D:200:PHE:HA	1:D:241:PHE:O	1.91	0.70
1:A:93:LEU:H	1:A:109:GLN:HE22	1.38	0.70
1:C:103:ARG:HB2	1:D:119:LEU:HD21	1.72	0.70
1:E:32:ILE:HB	1:E:53:VAL:HG12	1.73	0.70
1:F:42:ALA:HB1	1:F:45:LEU:HB3	1.73	0.70
1:B:115:THR:HG21	2:B:277:HOH:O	1.92	0.70
1:C:233:ASP:H	1:C:236:HIS:CD2	2.09	0.70
1:F:84:LEU:HD23	1:F:84:LEU:C	2.17	0.69
1:A:249:GLU:O	1:A:250:PHE:HB2	1.92	0.69
1:B:6:VAL:HG11	1:B:23:LEU:HD21	1.74	0.69
1:B:69:PHE:O	1:B:73:VAL:HG12	1.92	0.69
1:C:163:GLY:HA2	1:C:173:ARG:NE	2.07	0.69
1:F:158:THR:HG21	1:F:249:GLU:CB	2.23	0.69
1:C:164:SER:HB3	1:D:189:ASP:OD2	1.93	0.68
1:F:190:LEU:HD13	1:F:195:VAL:HG11	1.75	0.68
1:E:187:ALA:O	1:E:191:LYS:HE3	1.94	0.68
1:A:93:LEU:H	1:A:109:GLN:NE2	1.92	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:73:VAL:O	1:D:77:VAL:HG22	1.94	0.68
1:D:221:SER:HB3	1:D:243:ARG:HH11	1.58	0.68
1:B:198:VAL:C	1:B:199:ASN:HD22	2.01	0.68
1:B:6:VAL:HG13	1:B:84:LEU:HD13	1.76	0.67
1:A:111:ASP:HA	1:A:115:THR:CG2	2.24	0.67
1:C:138:ASP:HA	1:C:194:ASN:ND2	2.08	0.67
1:C:91:VAL:HG12	1:C:112:VAL:HG11	1.75	0.67
1:E:19:LEU:HD13	1:E:86:ILE:HD13	1.76	0.67
1:F:33:ILE:HD13	1:F:73:VAL:HG12	1.77	0.67
1:A:198:VAL:C	1:A:199:ASN:HD22	2.03	0.67
1:E:187:ALA:HB2	1:E:197:VAL:HG22	1.76	0.67
1:B:139:GLN:HB2	2:B:251:HOH:O	1.94	0.66
1:C:13:ARG:HH22	1:C:88:ASN:HD21	1.40	0.66
1:E:164:SER:HA	1:E:167:PHE:O	1.95	0.66
1:F:66:LEU:O	1:F:70:VAL:HG23	1.94	0.66
1:F:187:ALA:HA	1:F:197:VAL:CG1	2.25	0.66
1:A:173:ARG:HH22	1:C:166:GLN:NE2	1.94	0.66
1:B:242:MET:SD	1:B:246:LYS:HD3	2.36	0.66
1:D:249:GLU:O	1:D:250:PHE:HB2	1.96	0.66
1:F:70:VAL:O	1:F:73:VAL:HG22	1.96	0.66
1:A:103:ARG:HB2	1:B:119:LEU:HD21	1.77	0.65
1:B:249:GLU:O	1:B:250:PHE:HB2	1.94	0.65
1:C:19:LEU:HD13	1:C:86:ILE:HD13	1.78	0.65
1:D:162:SER:O	1:D:163:GLY:O	2.14	0.65
1:E:228:SER:HB3	1:E:245:LEU:HG	1.78	0.65
1:A:7:VAL:O	1:A:86:ILE:HG22	1.97	0.65
1:F:93:LEU:H	1:F:109:GLN:NE2	1.95	0.65
1:B:233:ASP:H	1:B:236:HIS:CD2	2.11	0.65
1:D:187:ALA:HB2	1:D:197:VAL:CG2	2.27	0.64
1:A:50:ASP:OD1	1:A:50:ASP:C	2.39	0.64
1:D:93:LEU:HD23	1:D:109:GLN:CD	2.22	0.64
1:B:160:ASN:ND2	1:B:163:GLY:H	1.94	0.64
1:F:190:LEU:HD13	1:F:195:VAL:CG1	2.27	0.64
1:D:15:ILE:H	1:D:15:ILE:CD1	2.10	0.64
1:F:18:GLY:HA3	1:F:222:THR:HG21	1.78	0.64
1:E:38:ASP:OD2	1:E:41:LYS:HB2	1.97	0.64
1:F:73:VAL:O	1:F:77:VAL:HG22	1.96	0.64
1:C:250:PHE:HD2	1:D:158:THR:HG22	1.64	0.63
1:D:232:LEU:HA	1:D:236:HIS:HD2	1.63	0.63
1:E:86:ILE:HA	1:E:148:ILE:O	1.99	0.63
1:F:6:VAL:HG11	1:F:23:LEU:HD21	1.81	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:233:ASP:H	1:F:236:HIS:CD2	2.16	0.63
1:E:52:ARG:HG3	1:E:52:ARG:HH11	1.64	0.62
1:F:127:LEU:H	1:F:127:LEU:CD1	2.11	0.62
1:F:158:THR:HG21	1:F:249:GLU:HB2	1.80	0.62
1:D:30:ARG:HG3	1:D:30:ARG:HH11	1.64	0.62
1:A:185:THR:HG23	1:B:163:GLY:HA2	1.81	0.62
1:F:18:GLY:O	1:F:22:GLN:HG2	2.00	0.62
1:D:69:PHE:O	1:D:73:VAL:HG23	1.98	0.62
1:F:63:ASP:O	1:F:66:LEU:HB2	2.00	0.62
1:A:86:ILE:C	1:A:86:ILE:HD13	2.25	0.61
1:C:138:ASP:HA	1:C:194:ASN:HD21	1.65	0.61
1:C:13:ARG:HH22	1:C:88:ASN:ND2	1.99	0.61
1:B:160:ASN:HD21	1:B:163:GLY:N	1.97	0.61
1:C:13:ARG:HE	1:C:13:ARG:C	2.07	0.61
1:D:233:ASP:H	1:D:236:HIS:CD2	2.16	0.61
1:B:92:LEU:C	1:B:92:LEU:HD23	2.26	0.61
1:D:242:MET:HB3	1:D:248:TYR:CE2	2.36	0.61
1:F:158:THR:HG22	1:F:250:PHE:O	2.00	0.61
1:D:125:LEU:N	1:D:126:PRO:HD2	2.16	0.61
1:E:233:ASP:H	1:E:236:HIS:HD2	1.48	0.61
1:D:15:ILE:HD12	1:D:15:ILE:N	2.11	0.60
1:F:52:ARG:HG2	1:F:52:ARG:HH11	1.65	0.60
1:F:63:ASP:HA	1:F:66:LEU:HD12	1.83	0.60
1:C:19:LEU:HD13	1:C:86:ILE:CD1	2.31	0.60
1:E:160:ASN:HD21	1:E:163:GLY:HA3	1.66	0.60
1:B:220:GLN:CD	1:B:220:GLN:H	2.09	0.60
1:C:19:LEU:O	1:C:23:LEU:HD22	2.01	0.60
1:A:26:ASP:O	1:A:52:ARG:NH2	2.34	0.60
1:B:190:LEU:HB3	1:B:195:VAL:HG13	1.83	0.60
1:F:83:SER:O	1:F:145:ALA:HA	2.01	0.60
1:C:37:ARG:HA	1:C:58:LEU:O	2.02	0.60
1:C:69:PHE:O	1:C:73:VAL:HG12	2.02	0.60
1:A:135:GLU:O	1:A:135:GLU:HG3	2.01	0.60
1:D:152:SER:HB3	1:D:176:LYS:HG3	1.83	0.60
1:A:8:VAL:HG22	1:A:86:ILE:CG2	2.32	0.59
1:E:42:ALA:O	1:E:46:LYS:HG3	2.02	0.59
1:E:21:GLN:O	1:E:24:VAL:HG22	2.02	0.59
1:A:233:ASP:H	1:A:236:HIS:HD2	1.49	0.59
1:C:50:ASP:HB3	1:C:53:VAL:HG13	1.84	0.59
1:C:17:LEU:O	1:C:21:GLN:HG3	2.03	0.59
1:C:44:GLU:O	1:C:47:SER:HB3	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:3:PRO:HB3	1:E:142:VAL:O	2.02	0.59
1:E:20:VAL:HA	1:E:23:LEU:HD23	1.85	0.59
1:E:38:ASP:OD1	1:E:41:LYS:HD3	2.02	0.59
1:C:71:SER:O	1:C:75:GLU:HG2	2.02	0.58
1:D:243:ARG:O	1:D:244:ASN:CG	2.45	0.58
1:B:86:ILE:HA	1:B:148:ILE:O	2.04	0.58
1:A:158:THR:CG2	1:A:250:PHE:OXT	2.52	0.58
1:D:19:LEU:HD13	1:D:86:ILE:HD11	1.86	0.58
1:C:218:VAL:HG23	1:C:220:GLN:H	1.68	0.58
1:E:232:LEU:HA	1:E:236:HIS:CD2	2.38	0.58
1:F:126:PRO:O	1:F:130:ASN:HB2	2.03	0.58
1:A:190:LEU:HB3	1:A:195:VAL:HG13	1.86	0.58
1:B:160:ASN:C	1:B:160:ASN:HD22	2.12	0.58
1:D:18:GLY:O	1:D:22:GLN:HG2	2.03	0.58
1:D:93:LEU:H	1:D:109:GLN:NE2	2.02	0.58
1:F:140:LEU:O	1:F:141:SER:O	2.22	0.58
1:A:244:ASN:O	1:A:245:LEU:HB2	2.02	0.58
1:B:233:ASP:N	1:B:236:HIS:HD2	1.97	0.58
1:A:8:VAL:HG22	1:A:86:ILE:HG21	1.84	0.57
1:E:97:THR:HG22	1:E:170:LEU:HD22	1.85	0.57
1:A:27:LYS:HA	1:A:52:ARG:NH2	2.20	0.57
1:A:50:ASP:O	1:A:53:VAL:HG12	2.03	0.57
1:C:198:VAL:C	1:C:199:ASN:HD22	2.12	0.57
1:F:42:ALA:O	1:F:44:GLU:N	2.36	0.57
1:A:134:LYS:O	1:A:135:GLU:HB3	2.04	0.57
1:A:160:ASN:ND2	1:A:163:GLY:H	2.02	0.57
1:B:86:ILE:HD13	1:B:86:ILE:C	2.28	0.57
1:E:12:ASN:O	1:E:13:ARG:HD3	2.04	0.57
1:E:186:LEU:HG	1:E:190:LEU:HD12	1.87	0.57
1:A:222:THR:O	1:A:226:ILE:HG12	2.04	0.57
1:E:17:LEU:O	1:E:21:GLN:HG3	2.05	0.57
1:C:92:LEU:CD2	1:C:92:LEU:C	2.78	0.56
1:B:93:LEU:N	1:B:109:GLN:HE22	2.00	0.56
1:B:190:LEU:HD13	1:B:195:VAL:HG11	1.87	0.56
1:F:239:ARG:HB2	1:F:241:PHE:CE2	2.41	0.56
1:A:189:ASP:OD2	1:B:164:SER:HB2	2.05	0.56
1:A:140:LEU:HB3	1:A:196:LEU:HD13	1.88	0.56
1:A:86:ILE:HD11	1:A:150:ILE:CG1	2.36	0.56
1:C:199:ASN:HD22	1:C:199:ASN:N	2.02	0.56
1:F:226:ILE:HD12	1:F:229:PHE:HB2	1.88	0.56
1:F:244:ASN:O	1:F:245:LEU:HB2	2.04	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:199:ASN:HD22	1:A:199:ASN:N	2.03	0.56
1:E:93:LEU:HD21	1:E:108:GLU:HB3	1.88	0.56
1:F:198:VAL:C	1:F:199:ASN:HD22	2.14	0.56
1:B:199:ASN:HD22	1:B:199:ASN:N	2.03	0.55
1:E:73:VAL:O	1:E:77:VAL:HG22	2.06	0.55
1:E:50:ASP:O	1:E:53:VAL:HG22	2.06	0.55
1:E:86:ILE:HD11	1:E:150:ILE:HD11	1.88	0.55
1:F:190:LEU:HB3	1:F:195:VAL:CG1	2.35	0.55
1:D:59:THR:HG1	1:D:62:CYS:HB3	1.71	0.55
1:E:158:THR:HG22	1:F:250:PHE:CD2	2.37	0.55
1:E:164:SER:HB3	1:F:189:ASP:OD2	2.06	0.55
1:D:93:LEU:N	1:D:93:LEU:HD22	2.22	0.55
1:F:7:VAL:HG22	1:F:33:ILE:HB	1.88	0.55
1:F:35:THR:HA	1:F:56:LEU:O	2.05	0.55
1:B:190:LEU:HB3	1:B:195:VAL:CG1	2.36	0.55
1:C:199:ASN:ND2	2:C:256:HOH:O	2.39	0.55
1:E:7:VAL:O	1:E:86:ILE:HG22	2.06	0.55
1:C:163:GLY:H	1:C:173:ARG:NH2	2.04	0.55
1:E:199:ASN:N	1:E:199:ASN:HD22	2.05	0.55
1:C:12:ASN:ND2	1:C:13:ARG:HG3	2.21	0.54
1:D:220:GLN:HA	1:D:220:GLN:OE1	2.05	0.54
1:D:22:GLN:HE21	1:D:22:GLN:HA	1.73	0.54
1:F:42:ALA:HB3	1:F:46:LYS:HG2	1.90	0.54
1:F:93:LEU:H	1:F:109:GLN:HE22	1.54	0.54
1:C:111:ASP:HA	1:C:115:THR:HG23	1.88	0.54
1:A:1:MET:HA	1:A:1:MET:CE	2.37	0.54
1:A:49:LYS:HE3	1:A:49:LYS:HA	1.89	0.54
1:C:91:VAL:HG22	1:C:92:LEU:N	2.22	0.54
1:C:163:GLY:H	1:C:173:ARG:HH21	1.55	0.54
1:D:72:LYS:O	1:D:76:ILE:HG12	2.08	0.54
1:F:84:LEU:CD2	1:F:86:ILE:HG13	2.38	0.54
1:F:160:ASN:HD21	1:F:163:GLY:H	1.55	0.54
1:B:244:ASN:O	1:B:245:LEU:HB2	2.07	0.54
1:D:18:GLY:HA3	1:D:222:THR:HG21	1.89	0.53
1:B:232:LEU:HA	1:B:236:HIS:CD2	2.43	0.53
1:D:190:LEU:HD13	1:D:195:VAL:HG11	1.90	0.53
1:F:122:GLN:HA	1:F:125:LEU:HG	1.90	0.53
1:E:158:THR:HG23	1:E:250:PHE:O	2.07	0.53
1:A:160:ASN:C	1:A:160:ASN:HD22	2.15	0.53
1:A:186:LEU:HD11	1:A:190:LEU:HD11	1.91	0.53
1:C:189:ASP:OD2	1:D:97:THR:HG23	2.07	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:92:LEU:C	1:E:92:LEU:HD23	2.33	0.53
1:F:129:LYS:HD2	1:F:193:ASP:OD2	2.09	0.53
1:E:69:PHE:O	1:E:73:VAL:HG13	2.08	0.53
1:B:39:VAL:O	1:B:46:LYS:HD2	2.08	0.53
1:C:86:ILE:HD12	1:C:150:ILE:HD11	1.90	0.53
1:B:59:THR:HA	2:B:279:HOH:O	2.09	0.53
1:D:224:GLU:CD	1:D:243:ARG:HD2	2.34	0.53
1:F:249:GLU:O	1:F:250:PHE:HB2	2.07	0.53
1:C:39:VAL:HG22	1:C:39:VAL:O	2.09	0.53
1:A:13:ARG:HB2	2:A:264:HOH:O	2.09	0.52
1:D:57:PRO:O	1:D:58:LEU:HB2	2.09	0.52
1:A:86:ILE:HG23	1:A:86:ILE:O	2.10	0.52
1:C:13:ARG:NH2	1:C:88:ASN:ND2	2.55	0.52
1:F:176:LYS:O	1:F:179:ILE:HG22	2.09	0.52
1:D:12:ASN:HA	1:D:45:LEU:HD12	1.91	0.52
1:C:228:SER:O	1:C:232:LEU:HG	2.08	0.52
1:D:45:LEU:HD22	1:D:55:VAL:HG22	1.91	0.52
1:B:239:ARG:HB2	1:B:241:PHE:CZ	2.45	0.52
1:C:122:GLN:HE21	1:D:97:THR:HA	1.74	0.52
1:D:92:LEU:HD23	1:D:93:LEU:O	2.10	0.52
1:D:7:VAL:HG22	1:D:33:ILE:HB	1.92	0.52
1:A:86:ILE:HD11	1:A:150:ILE:HG12	1.91	0.52
1:F:97:THR:HG22	1:F:170:LEU:HD22	1.91	0.52
1:B:222:THR:O	1:B:226:ILE:HG12	2.10	0.52
1:F:62:CYS:O	1:F:66:LEU:HG	2.10	0.52
1:F:233:ASP:N	1:F:236:HIS:HD2	2.06	0.52
1:C:32:ILE:O	1:C:53:VAL:HA	2.10	0.51
1:D:92:LEU:HD23	1:D:92:LEU:C	2.35	0.51
1:E:202:PRO:HG2	1:E:203:GLY:H	1.75	0.51
1:F:15:ILE:N	1:F:15:ILE:HD12	2.25	0.51
1:B:222:THR:HA	1:B:225:LEU:HD12	1.91	0.51
1:C:244:ASN:O	1:C:245:LEU:HB2	2.11	0.51
1:E:103:ARG:NH1	1:F:111:ASP:OD1	2.43	0.51
1:E:93:LEU:H	1:E:109:GLN:NE2	2.08	0.51
1:D:160:ASN:C	1:D:160:ASN:HD22	2.18	0.51
1:B:86:ILE:HD13	1:B:86:ILE:O	2.11	0.51
1:C:20:VAL:O	1:C:24:VAL:HG13	2.11	0.51
1:E:39:VAL:HG13	1:E:40:GLU:N	2.26	0.51
1:F:224:GLU:O	1:F:227:SER:HB3	2.11	0.51
1:E:113:ASN:HB2	1:E:175:SER:HB2	1.92	0.51
1:F:43:THR:HA	1:F:46:LYS:HB2	1.92	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:132:ALA:HB2	1:C:145:ALA:CB	2.41	0.51
1:D:8:VAL:HA	1:D:86:ILE:HG23	1.93	0.51
1:D:221:SER:HB3	1:D:243:ARG:NH1	2.25	0.50
1:F:140:LEU:HD22	1:F:234:ASN:HA	1.93	0.50
1:B:93:LEU:CD1	1:B:108:GLU:HB3	2.41	0.50
1:F:69:PHE:O	1:F:73:VAL:HG13	2.12	0.50
1:F:115:THR:O	1:F:119:LEU:HD22	2.12	0.50
1:C:185:THR:HG21	1:D:170:LEU:HD13	1.93	0.50
1:C:190:LEU:HB3	1:C:195:VAL:HG13	1.94	0.50
1:F:50:ASP:OD2	1:F:52:ARG:NH1	2.43	0.50
1:F:120:LEU:O	1:F:124:LEU:HG	2.11	0.50
1:A:245:LEU:O	1:A:246:LYS:CB	2.59	0.50
1:C:186:LEU:HD11	1:C:190:LEU:HD11	1.93	0.50
1:D:50:ASP:O	1:D:53:VAL:HG12	2.12	0.50
1:E:40:GLU:C	1:E:41:LYS:HD2	2.37	0.50
1:B:93:LEU:HD11	1:B:108:GLU:HB3	1.92	0.50
1:C:188:VAL:HG11	1:D:163:GLY:H	1.77	0.50
1:D:222:THR:O	1:D:226:ILE:HG12	2.12	0.50
1:D:19:LEU:HD13	1:D:86:ILE:CD1	2.42	0.50
1:B:249:GLU:O	1:B:250:PHE:CB	2.60	0.49
1:A:72:LYS:O	1:A:76:ILE:HG13	2.11	0.49
1:D:30:ARG:HG3	1:D:30:ARG:NH1	2.27	0.49
1:A:204:TRP:CD1	1:C:168:PRO:HD3	2.48	0.49
1:C:86:ILE:CD1	1:C:150:ILE:HD11	2.43	0.49
1:F:5:SER:OG	1:F:83:SER:HB3	2.12	0.49
1:F:222:THR:HA	1:F:225:LEU:HD12	1.93	0.49
1:D:35:THR:HA	1:D:56:LEU:O	2.13	0.49
1:D:193:ASP:O	1:D:194:ASN:HB2	2.12	0.49
1:F:140:LEU:HD13	1:F:234:ASN:ND2	2.27	0.49
1:A:35:THR:HA	1:A:56:LEU:O	2.11	0.49
1:B:45:LEU:HD23	1:B:48:ILE:HD12	1.94	0.49
1:C:21:GLN:O	1:C:24:VAL:HG22	2.13	0.49
1:D:11:ALA:HA	1:D:16:GLY:HA3	1.95	0.49
1:D:140:LEU:O	1:D:233:ASP:HA	2.13	0.49
1:F:21:GLN:O	1:F:24:VAL:HG22	2.13	0.49
1:A:17:LEU:O	1:A:21:GLN:HG3	2.13	0.49
1:B:118:VAL:O	1:B:122:GLN:HG3	2.13	0.48
1:D:187:ALA:HB2	1:D:197:VAL:HG21	1.95	0.48
1:A:49:LYS:O	1:A:50:ASP:O	2.31	0.48
1:A:160:ASN:ND2	1:A:160:ASN:C	2.71	0.48
1:B:160:ASN:ND2	1:B:160:ASN:C	2.71	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:160:ASN:C	1:C:160:ASN:HD22	2.21	0.48
1:E:160:ASN:HD21	1:E:163:GLY:CA	2.25	0.48
1:D:233:ASP:N	1:D:236:HIS:HD2	2.08	0.48
1:E:92:LEU:HD23	1:E:92:LEU:O	2.14	0.48
1:F:125:LEU:HB2	1:F:126:PRO:HD3	1.94	0.48
1:F:187:ALA:CA	1:F:197:VAL:HG11	2.42	0.48
1:C:227:SER:O	1:C:231:LYS:HG3	2.13	0.48
1:D:59:THR:OG1	1:D:62:CYS:HB3	2.13	0.48
1:D:80:ASP:HA	1:D:134:LYS:HE3	1.94	0.48
1:A:242:MET:O	1:A:245:LEU:N	2.45	0.48
1:A:249:GLU:O	1:A:250:PHE:CB	2.60	0.48
1:E:50:ASP:OD1	1:E:52:ARG:HG3	2.14	0.48
1:E:92:LEU:HA	1:E:109:GLN:HE22	1.78	0.48
1:F:113:ASN:ND2	1:F:176:LYS:HE2	2.28	0.48
1:D:190:LEU:HB3	1:D:195:VAL:CG1	2.42	0.48
1:E:11:ALA:HB3	1:E:34:ALA:HB1	1.94	0.48
1:C:82:LEU:HD12	1:C:128:LEU:HD21	1.96	0.48
1:C:84:LEU:HD22	1:C:86:ILE:HG22	1.94	0.48
1:A:185:THR:HG23	1:B:163:GLY:CA	2.44	0.48
1:D:86:ILE:HD12	1:D:229:PHE:HE2	1.78	0.48
1:B:20:VAL:O	1:B:24:VAL:HG13	2.14	0.48
1:F:73:VAL:HG23	1:F:74:GLY:N	2.29	0.47
1:A:15:ILE:HD12	1:A:15:ILE:N	2.23	0.47
1:D:64:LYS:HA	1:D:64:LYS:CE	2.40	0.47
1:E:83:SER:O	1:E:145:ALA:HA	2.13	0.47
1:F:84:LEU:HD21	1:F:86:ILE:HG13	1.94	0.47
1:F:232:LEU:HA	1:F:236:HIS:CD2	2.48	0.47
1:E:185:THR:HG21	1:F:170:LEU:HD13	1.95	0.47
1:C:249:GLU:O	1:C:250:PHE:HB2	2.13	0.47
1:E:50:ASP:OD2	1:E:52:ARG:NH1	2.46	0.47
1:E:87:ASN:HB2	1:E:149:THR:HG23	1.96	0.47
1:C:86:ILE:HA	1:C:148:ILE:O	2.15	0.47
1:D:83:SER:O	1:D:145:ALA:HA	2.15	0.47
1:A:15:ILE:H	1:A:15:ILE:CD1	2.22	0.47
1:A:70:VAL:O	1:A:73:VAL:HG22	2.15	0.47
1:A:83:SER:O	1:A:145:ALA:HA	2.15	0.47
1:D:22:GLN:HA	1:D:22:GLN:NE2	2.29	0.47
1:F:196:LEU:HD21	1:F:232:LEU:HB3	1.96	0.47
1:A:86:ILE:HA	1:A:148:ILE:O	2.15	0.47
1:A:113:ASN:ND2	1:A:176:LYS:HE2	2.30	0.47
1:A:115:THR:HB	1:B:106:ILE:HG21	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:198:VAL:C	1:D:199:ASN:HD22	2.22	0.47
1:F:91:VAL:HG13	1:F:93:LEU:CD2	2.44	0.47
1:D:187:ALA:O	1:D:191:LYS:HE3	2.15	0.47
1:B:190:LEU:HD13	1:B:195:VAL:CG1	2.44	0.47
1:A:22:GLN:HE21	1:A:22:GLN:HA	1.80	0.46
1:D:49:LYS:O	1:D:50:ASP:C	2.58	0.46
1:E:196:LEU:HD12	1:E:237:ASN:OD1	2.15	0.46
1:D:45:LEU:O	1:D:45:LEU:HD23	2.16	0.46
1:F:158:THR:HG21	1:F:249:GLU:HB3	1.96	0.46
1:A:140:LEU:O	1:A:233:ASP:HA	2.15	0.46
1:D:188:VAL:O	1:D:191:LYS:HB2	2.16	0.46
1:C:102:ASN:OD1	1:C:104:ALA:HB3	2.16	0.46
1:C:31:HIS:HD2	2:C:265:HOH:O	1.97	0.46
1:C:38:ASP:C	1:C:40:GLU:H	2.24	0.46
1:C:73:VAL:O	1:C:77:VAL:HG22	2.16	0.46
1:E:125:LEU:N	1:E:126:PRO:CD	2.79	0.46
1:A:99:THR:O	1:A:100:GLU:C	2.58	0.46
1:C:164:SER:HB3	1:D:189:ASP:CG	2.41	0.46
1:B:9:THR:HA	1:B:35:THR:OG1	2.16	0.46
1:D:23:LEU:HB3	1:D:32:ILE:HD13	1.96	0.46
1:D:66:LEU:CD2	1:D:120:LEU:HD13	2.45	0.46
1:F:123:LYS:O	1:F:124:LEU:HD23	2.15	0.46
1:A:50:ASP:OD2	1:A:52:ARG:NH1	2.48	0.46
1:C:12:ASN:CG	1:C:13:ARG:HG3	2.40	0.46
1:E:113:ASN:ND2	1:E:176:LYS:HE2	2.31	0.46
1:C:1:MET:HG3	1:C:2:SER:H	1.81	0.45
1:F:199:ASN:HD22	1:F:199:ASN:N	2.11	0.45
1:B:22:GLN:OE1	1:B:223:ALA:HB2	2.16	0.45
1:C:189:ASP:OD2	1:D:97:THR:CG2	2.64	0.45
1:E:162:SER:O	1:F:188:VAL:HG11	2.16	0.45
1:E:91:VAL:HG22	1:E:92:LEU:N	2.31	0.45
1:B:19:LEU:O	1:B:23:LEU:HB2	2.17	0.45
1:B:170:LEU:O	1:B:174:MET:HG3	2.17	0.45
1:D:249:GLU:O	1:D:250:PHE:CB	2.64	0.45
1:F:140:LEU:HD23	1:F:196:LEU:HD13	1.99	0.45
1:A:22:GLN:HB3	1:A:226:ILE:HG13	1.97	0.45
1:C:160:ASN:C	1:C:160:ASN:ND2	2.73	0.45
1:D:17:LEU:O	1:D:21:GLN:HG3	2.16	0.45
1:B:92:LEU:C	1:B:92:LEU:CD2	2.90	0.45
1:D:37:ARG:HG3	1:D:59:THR:HG22	1.99	0.45
1:E:11:ALA:O	1:E:17:LEU:HB2	2.14	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:66:LEU:O	1:A:70:VAL:HG23	2.17	0.45
1:E:12:ASN:HD22	1:E:13:ARG:HG2	1.78	0.45
1:E:242:MET:HG2	1:E:248:TYR:CE2	2.51	0.45
1:F:228:SER:HB2	1:F:245:LEU:HG	1.99	0.45
1:A:79:SER:O	1:A:134:LYS:HE2	2.16	0.45
1:E:178:ALA:O	1:E:181:MET:HB3	2.17	0.45
1:A:10:GLY:O	1:A:16:GLY:HA3	2.16	0.45
1:C:92:LEU:HD23	1:C:93:LEU:N	2.31	0.45
1:E:84:LEU:CD2	1:E:85:LEU:N	2.80	0.45
1:C:164:SER:HA	1:C:167:PHE:O	2.17	0.44
1:E:6:VAL:HG12	1:E:7:VAL:N	2.31	0.44
1:E:12:ASN:HD22	1:E:13:ARG:NE	2.15	0.44
1:F:68:THR:HG22	1:F:72:LYS:HE3	1.99	0.44
1:F:113:ASN:HD22	1:F:176:LYS:HE2	1.82	0.44
1:C:222:THR:O	1:C:226:ILE:HG12	2.17	0.44
1:A:39:VAL:C	1:A:41:LYS:N	2.76	0.44
1:B:49:LYS:HB2	1:B:49:LYS:NZ	2.32	0.44
1:C:198:VAL:HG22	1:C:236:HIS:O	2.18	0.44
1:D:176:LYS:O	1:D:179:ILE:HG22	2.17	0.44
1:D:242:MET:O	1:D:243:ARG:O	2.35	0.44
1:E:19:LEU:HD13	1:E:86:ILE:CD1	2.45	0.44
1:E:60:VAL:HB	1:E:116:SER:HB3	1.98	0.44
1:D:242:MET:C	1:D:243:ARG:O	2.58	0.44
1:E:184:ARG:HG2	1:F:160:ASN:HB3	2.00	0.44
1:F:17:LEU:O	1:F:21:GLN:HG3	2.18	0.44
1:A:91:VAL:HG22	1:A:92:LEU:N	2.33	0.44
1:B:187:ALA:HA	1:B:197:VAL:HG13	2.00	0.44
1:C:163:GLY:N	1:C:173:ARG:HH21	2.13	0.44
1:C:222:THR:HA	1:C:225:LEU:HD12	1.99	0.44
1:D:57:PRO:O	1:D:58:LEU:CB	2.66	0.44
1:D:7:VAL:O	1:D:86:ILE:HG23	2.17	0.44
1:D:92:LEU:HA	1:D:109:GLN:HE22	1.83	0.44
1:E:139:GLN:O	1:E:139:GLN:HG2	2.18	0.44
1:B:85:LEU:HD22	1:B:124:LEU:HD12	2.00	0.44
1:A:242:MET:HB2	1:A:245:LEU:O	2.18	0.44
1:C:239:ARG:HB2	1:C:241:PHE:CZ	2.53	0.44
1:C:122:GLN:HA	1:C:125:LEU:HG	2.00	0.43
1:D:45:LEU:HD22	1:D:55:VAL:CG2	2.48	0.43
1:D:227:SER:O	1:D:231:LYS:HG3	2.18	0.43
1:E:233:ASP:N	1:E:236:HIS:HD2	2.16	0.43
1:D:87:ASN:CG	1:D:117:VAL:HG13	2.43	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:9:THR:HB	1:F:89:ALA:HB2	1.99	0.43
1:F:30:ARG:O	1:F:52:ARG:HB2	2.18	0.43
1:C:207:THR:O	1:C:207:THR:CG2	2.66	0.43
1:D:233:ASP:OD1	1:D:233:ASP:C	2.62	0.43
1:E:113:ASN:O	1:E:179:ILE:HG13	2.19	0.43
1:F:160:ASN:C	1:F:160:ASN:ND2	2.76	0.43
1:A:9:THR:O	1:A:88:ASN:HB3	2.18	0.43
1:A:122:GLN:HB3	1:B:101:PRO:HG3	2.00	0.43
1:B:1:MET:HA	1:B:1:MET:CE	2.42	0.43
1:B:146:ALA:HA	1:B:196:LEU:O	2.18	0.43
1:C:163:GLY:O	1:D:189:ASP:OD1	2.36	0.43
1:F:151:SER:HB3	1:F:240:PHE:CE1	2.53	0.43
1:F:169:VAL:O	1:F:173:ARG:HG3	2.19	0.43
1:B:178:ALA:O	1:B:181:MET:HB3	2.18	0.43
1:C:1:MET:HG3	1:C:2:SER:N	2.34	0.43
1:E:52:ARG:HG3	1:E:52:ARG:NH1	2.31	0.43
1:A:49:LYS:O	1:A:50:ASP:C	2.62	0.43
1:B:160:ASN:HD22	1:B:161:THR:N	2.17	0.43
1:D:43:THR:HA	1:D:46:LYS:HD3	2.01	0.43
1:D:45:LEU:C	1:D:47:SER:H	2.27	0.43
1:E:70:VAL:HG21	1:E:123:LYS:HB3	2.00	0.43
1:E:140:LEU:HB3	1:E:196:LEU:HB2	1.99	0.43
1:A:22:GLN:HA	1:A:22:GLN:NE2	2.33	0.43
1:A:69:PHE:O	1:A:73:VAL:HG13	2.19	0.43
1:C:162:SER:OG	1:C:163:GLY:N	2.52	0.43
1:D:88:ASN:HA	1:D:150:ILE:HD13	2.01	0.43
1:A:18:GLY:O	1:A:22:GLN:HG2	2.18	0.43
1:B:70:VAL:HA	1:B:73:VAL:CG1	2.48	0.43
1:C:207:THR:O	1:C:207:THR:HG22	2.18	0.43
1:D:86:ILE:HD12	1:D:229:PHE:CE2	2.54	0.43
1:D:91:VAL:HG13	1:D:93:LEU:HD22	2.00	0.43
1:B:50:ASP:OD2	1:B:52:ARG:NH1	2.52	0.43
1:F:202:PRO:C	1:F:204:TRP:H	2.26	0.43
1:C:125:LEU:N	1:C:126:PRO:CD	2.82	0.42
1:D:8:VAL:HG22	1:D:86:ILE:HG21	2.01	0.42
1:E:26:ASP:C	1:E:28:ASN:H	2.27	0.42
1:E:157:ILE:O	1:E:160:ASN:HB3	2.18	0.42
1:F:20:VAL:O	1:F:24:VAL:HG13	2.19	0.42
1:B:157:ILE:N	1:B:250:PHE:O	2.40	0.42
1:D:9:THR:O	1:D:88:ASN:HB3	2.19	0.42
1:D:97:THR:HG23	1:D:164:SER:OG	2.19	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:15:ILE:HG12	1:F:202:PRO:HG3	2.00	0.42
1:F:200:PHE:HZ	1:F:232:LEU:HD21	1.85	0.42
1:A:50:ASP:HA	2:A:293:HOH:O	2.20	0.42
1:E:119:LEU:HD13	1:F:106:ILE:HD12	2.01	0.42
1:A:127:LEU:HG	2:A:262:HOH:O	2.20	0.42
1:D:87:ASN:O	1:D:150:ILE:HD13	2.19	0.42
1:F:160:ASN:C	1:F:160:ASN:HD22	2.26	0.42
1:A:100:GLU:O	1:A:101:PRO:C	2.61	0.42
1:B:187:ALA:HA	1:B:197:VAL:CG1	2.49	0.42
1:C:49:LYS:HD3	1:C:49:LYS:N	2.34	0.42
1:C:92:LEU:C	1:C:92:LEU:HD22	2.44	0.42
1:B:64:LYS:HD3	1:B:64:LYS:HA	1.76	0.42
1:D:7:VAL:O	1:D:7:VAL:HG12	2.19	0.42
1:E:8:VAL:HG22	1:E:86:ILE:CG2	2.49	0.42
1:F:151:SER:HB3	1:F:240:PHE:HE1	1.84	0.42
1:F:231:LYS:HE2	1:F:231:LYS:CA	2.38	0.42
1:D:197:VAL:O	1:D:197:VAL:HG23	2.20	0.42
1:E:18:GLY:O	1:E:22:GLN:HG2	2.20	0.42
1:E:82:LEU:HD12	1:E:124:LEU:HD22	2.02	0.42
1:E:160:ASN:HD21	1:E:163:GLY:N	2.18	0.42
1:E:239:ARG:HB2	1:E:241:PHE:CE1	2.55	0.42
1:E:39:VAL:HG23	1:E:55:VAL:HG11	2.02	0.42
1:F:26:ASP:OD1	1:F:29:ILE:HG13	2.19	0.41
1:B:23:LEU:HD13	1:B:226:ILE:HD12	2.02	0.41
1:C:82:LEU:HD12	1:C:128:LEU:CD2	2.50	0.41
1:E:225:LEU:HD23	1:E:243:ARG:HA	2.03	0.41
1:F:93:LEU:N	1:F:93:LEU:HD22	2.35	0.41
1:F:140:LEU:HD23	1:F:196:LEU:CD1	2.49	0.41
1:B:132:ALA:HB2	1:B:145:ALA:CB	2.50	0.41
1:B:176:LYS:O	1:B:179:ILE:HG22	2.20	0.41
1:C:189:ASP:OD1	1:D:163:GLY:HA3	2.20	0.41
1:D:19:LEU:HD22	1:D:226:ILE:HD11	2.02	0.41
1:B:138:ASP:HA	1:B:194:ASN:ND2	2.35	0.41
1:C:190:LEU:HB3	1:C:195:VAL:CG1	2.50	0.41
1:B:23:LEU:HD12	1:B:23:LEU:HA	1.93	0.41
1:C:39:VAL:HG23	1:C:55:VAL:HG12	2.02	0.41
1:D:160:ASN:C	1:D:160:ASN:ND2	2.78	0.41
1:E:179:ILE:HD13	1:E:179:ILE:HA	1.88	0.41
1:A:233:ASP:H	1:A:236:HIS:CD2	2.34	0.41
1:C:50:ASP:OD1	1:C:52:ARG:HG3	2.20	0.41
1:F:50:ASP:OD1	1:F:52:ARG:HG2	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:32:ILE:O	1:D:53:VAL:HA	2.21	0.41
1:E:93:LEU:CD2	1:E:108:GLU:HB3	2.50	0.41
1:F:117:VAL:HG11	1:F:179:ILE:HD11	2.03	0.41
1:D:190:LEU:HD13	1:D:195:VAL:CG1	2.50	0.41
1:D:227:SER:O	1:D:231:LYS:HE3	2.21	0.41
1:A:39:VAL:O	1:A:41:LYS:N	2.53	0.41
1:B:9:THR:O	1:B:88:ASN:HB3	2.21	0.41
1:C:12:ASN:C	1:C:13:ARG:HG3	2.45	0.41
1:E:84:LEU:HD21	1:E:148:ILE:HD13	2.02	0.41
1:E:32:ILE:O	1:E:53:VAL:HA	2.20	0.41
1:F:242:MET:HG2	1:F:248:TYR:CE2	2.56	0.41
1:A:157:ILE:HG12	1:B:181:MET:HE2	2.02	0.40
1:B:144:ARG:HB3	2:B:289:HOH:O	2.21	0.40
1:D:64:LYS:HE3	1:D:64:LYS:CA	2.47	0.40
1:D:176:LYS:HA	1:D:176:LYS:HD3	1.87	0.40
1:F:221:SER:O	1:F:225:LEU:HG	2.20	0.40
1:A:49:LYS:HA	1:A:49:LYS:CE	2.52	0.40
1:A:135:GLU:O	1:A:135:GLU:CG	2.68	0.40
1:D:140:LEU:HB2	1:D:234:ASN:OD1	2.21	0.40
1:E:226:ILE:HA	1:E:229:PHE:CD2	2.55	0.40
1:F:109:GLN:HG2	1:F:171:ALA:CB	2.51	0.40
1:F:146:ALA:HA	1:F:196:LEU:O	2.22	0.40
1:D:151:SER:HA	1:D:176:LYS:HD2	2.03	0.40
1:E:198:VAL:C	1:E:199:ASN:HD22	2.30	0.40
1:F:52:ARG:HH11	1:F:52:ARG:CG	2.29	0.40
1:A:182:PHE:HD1	1:B:170:LEU:HD11	1.85	0.40
1:B:79:SER:O	1:B:130:ASN:ND2	2.53	0.40
1:F:91:VAL:HG13	1:F:93:LEU:HD21	2.03	0.40
1:D:52:ARG:HG2	1:D:52:ARG:H	1.77	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	233/250 (93%)	214 (92%)	13 (6%)	6 (3%)	4	7
1	B	231/250 (92%)	211 (91%)	18 (8%)	2 (1%)	14	30
1	C	236/250 (94%)	224 (95%)	10 (4%)	2 (1%)	16	34
1	D	232/250 (93%)	209 (90%)	15 (6%)	8 (3%)	3	4
1	E	231/250 (92%)	212 (92%)	16 (7%)	3 (1%)	9	21
1	F	220/250 (88%)	194 (88%)	23 (10%)	3 (1%)	9	19
All	All	1383/1500 (92%)	1264 (91%)	95 (7%)	24 (2%)	7	15

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	49	LYS
1	A	50	ASP
1	A	135	GLU
1	C	163	GLY
1	D	58	LEU
1	D	163	GLY
1	E	163	GLY
1	F	43	THR
1	F	141	SER
1	B	60	VAL
1	C	39	VAL
1	D	50	ASP
1	D	139	GLN
1	D	243	ARG
1	E	27	LYS
1	F	144	ARG
1	A	246	LYS
1	B	166	GLN
1	D	234	ASN
1	A	134	LYS
1	A	245	LEU
1	E	14	GLY
1	D	165	ALA
1	D	101	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	203/212 (96%)	181 (89%)	22 (11%)	6	13
1	B	201/212 (95%)	186 (92%)	15 (8%)	12	28
1	C	206/212 (97%)	187 (91%)	19 (9%)	8	19
1	D	202/212 (95%)	190 (94%)	12 (6%)	18	38
1	E	201/212 (95%)	191 (95%)	10 (5%)	22	46
1	F	198/212 (93%)	192 (97%)	6 (3%)	36	64
All	All	1211/1272 (95%)	1127 (93%)	84 (7%)	14	32

All (84) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	MET
1	A	2	SER
1	A	13	ARG
1	A	23	LEU
1	A	39	VAL
1	A	45	LEU
1	A	49	LYS
1	A	50	ASP
1	A	52	ARG
1	A	53	VAL
1	A	71	SER
1	A	82	LEU
1	A	84	LEU
1	A	86	ILE
1	A	93	LEU
1	A	115	THR
1	A	119	LEU
1	A	158	THR
1	A	179	ILE
1	A	195	VAL
1	A	197	VAL
1	A	199	ASN
1	B	23	LEU
1	B	49	LYS
1	B	52	ARG
1	B	59	THR
1	B	73	VAL

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Mol	Chain	Res	Type
1	B	82	LEU
1	B	84	LEU
1	B	86	ILE
1	B	115	THR
1	B	119	LEU
1	B	140	LEU
1	B	191	LYS
1	B	195	VAL
1	B	197	VAL
1	B	199	ASN
1	C	13	ARG
1	C	23	LEU
1	C	37	ARG
1	C	45	LEU
1	C	49	LYS
1	C	53	VAL
1	C	73	VAL
1	C	84	LEU
1	C	86	ILE
1	C	92	LEU
1	C	115	THR
1	C	119	LEU
1	C	138	ASP
1	C	140	LEU
1	C	158	THR
1	C	159	ASP
1	C	195	VAL
1	C	197	VAL
1	C	199	ASN
1	D	23	LEU
1	D	52	ARG
1	D	64	LYS
1	D	68	THR
1	D	84	LEU
1	D	93	LEU
1	D	115	THR
1	D	119	LEU
1	D	139	GLN
1	D	140	LEU
1	D	191	LYS
1	D	195	VAL
1	E	23	LEU

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Mol	Chain	Res	Type
1	E	59	THR
1	E	73	VAL
1	E	82	LEU
1	E	84	LEU
1	E	119	LEU
1	E	138	ASP
1	E	179	ILE
1	E	197	VAL
1	E	199	ASN
1	F	23	LEU
1	F	45	LEU
1	F	115	THR
1	F	119	LEU
1	F	197	VAL
1	F	204	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (59) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	22	GLN
1	A	109	GLN
1	A	130	ASN
1	A	139	GLN
1	A	160	ASN
1	A	166	GLN
1	A	199	ASN
1	A	236	HIS
1	B	22	GLN
1	B	31	HIS
1	B	54	HIS
1	B	87	ASN
1	B	88	ASN
1	B	109	GLN
1	B	160	ASN
1	B	194	ASN
1	B	199	ASN
1	B	230	ASN
1	B	236	HIS
1	C	21	GLN
1	C	22	GLN
1	C	31	HIS
1	C	87	ASN

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Mol	Chain	Res	Type
1	C	88	ASN
1	C	109	GLN
1	C	122	GLN
1	C	160	ASN
1	C	166	GLN
1	C	194	ASN
1	C	199	ASN
1	C	206	GLN
1	C	236	HIS
1	D	22	GLN
1	D	31	HIS
1	D	54	HIS
1	D	87	ASN
1	D	102	ASN
1	D	109	GLN
1	D	160	ASN
1	D	194	ASN
1	D	199	ASN
1	D	236	HIS
1	E	12	ASN
1	E	98	ASN
1	E	109	GLN
1	E	122	GLN
1	E	160	ASN
1	E	194	ASN
1	E	199	ASN
1	E	220	GLN
1	E	236	HIS
1	F	22	GLN
1	F	31	HIS
1	F	109	GLN
1	F	122	GLN
1	F	160	ASN
1	F	194	ASN
1	F	199	ASN
1	F	236	HIS

5.3.3 RNA ⓘ

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	237/250 (94%)	-0.25	1 (0%) 88 86	26, 41, 61, 74	0
1	B	235/250 (94%)	-0.23	1 (0%) 88 86	27, 44, 64, 73	0
1	C	240/250 (96%)	-0.31	2 (0%) 82 80	27, 38, 59, 71	0
1	D	236/250 (94%)	0.17	4 (1%) 69 64	29, 48, 79, 92	0
1	E	235/250 (94%)	-0.01	2 (0%) 81 78	33, 50, 69, 82	0
1	F	230/250 (92%)	0.33	9 (3%) 43 38	34, 58, 79, 88	0
All	All	1413/1500 (94%)	-0.05	19 (1%) 75 71	26, 45, 71, 92	0

All (19) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	218	VAL	4.3
1	F	204	TRP	3.7
1	C	218	VAL	3.3
1	F	202	PRO	2.8
1	F	195	VAL	2.8
1	E	218	VAL	2.7
1	D	5	SER	2.5
1	F	80	ASP	2.3
1	D	221	SER	2.3
1	F	140	LEU	2.2
1	F	83	SER	2.2
1	D	23	LEU	2.1
1	F	219	GLU	2.1
1	C	12	ASN	2.1
1	F	203	GLY	2.1
1	D	41	LYS	2.0
1	B	220	GLN	2.0
1	E	43	THR	2.0
1	F	77	VAL	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.