



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 02:26 AM UTC

PDB ID : 3ALE / pdb_00003ale
Title : A type III polyketide synthase that produces diarylheptanoid
Authors : Morita, H.; Kato, R.; Sugio, S.; Abe, I.
Deposited on : 2010-08-03
Resolution : 2.50 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

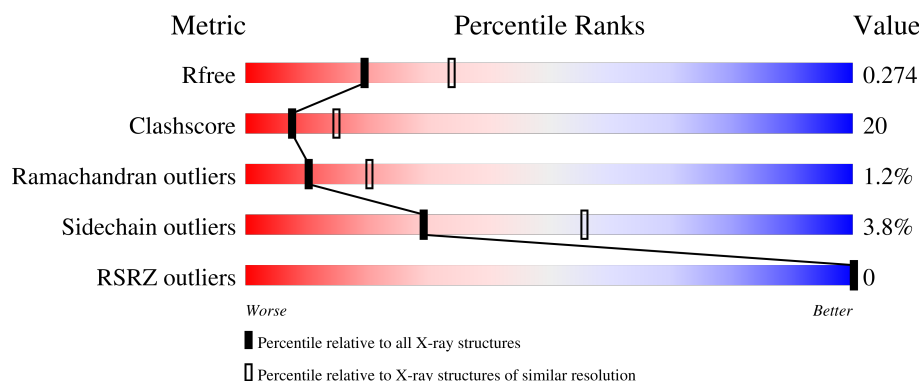
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	416	 60% 25% • 11%
1	B	416	 52% 33% • 12%
1	C	416	 52% 32% • 12%
1	D	416	 56% 29% • 13%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 11223 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 Os07g0271500 protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	369	Total	C	N	O	S	0	0	0
			2805	1764	504	519	18			
1	B	366	Total	C	N	O	S	0	0	0
			2778	1746	498	516	18			
1	C	364	Total	C	N	O	S	0	0	0
			2768	1741	496	513	18			
1	D	361	Total	C	N	O	S	0	0	0
			2747	1730	493	506	18			

There are 60 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-13	MET	-	expression tag	UNP Q8LIL0
A	-12	GLY	-	expression tag	UNP Q8LIL0
A	-11	SER	-	expression tag	UNP Q8LIL0
A	-10	SER	-	expression tag	UNP Q8LIL0
A	-9	HIS	-	expression tag	UNP Q8LIL0
A	-8	HIS	-	expression tag	UNP Q8LIL0
A	-7	HIS	-	expression tag	UNP Q8LIL0
A	-6	HIS	-	expression tag	UNP Q8LIL0
A	-5	HIS	-	expression tag	UNP Q8LIL0
A	-4	HIS	-	expression tag	UNP Q8LIL0
A	-3	SER	-	expression tag	UNP Q8LIL0
A	-2	GLN	-	expression tag	UNP Q8LIL0
A	-1	ASP	-	expression tag	UNP Q8LIL0
A	0	PRO	-	expression tag	UNP Q8LIL0
A	46	ILE	PHE	SEE REMARK 999	UNP Q8LIL0
B	-13	MET	-	expression tag	UNP Q8LIL0
B	-12	GLY	-	expression tag	UNP Q8LIL0
B	-11	SER	-	expression tag	UNP Q8LIL0
B	-10	SER	-	expression tag	UNP Q8LIL0
B	-9	HIS	-	expression tag	UNP Q8LIL0
B	-8	HIS	-	expression tag	UNP Q8LIL0

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-7	HIS	-	expression tag	UNP Q8LIL0
B	-6	HIS	-	expression tag	UNP Q8LIL0
B	-5	HIS	-	expression tag	UNP Q8LIL0
B	-4	HIS	-	expression tag	UNP Q8LIL0
B	-3	SER	-	expression tag	UNP Q8LIL0
B	-2	GLN	-	expression tag	UNP Q8LIL0
B	-1	ASP	-	expression tag	UNP Q8LIL0
B	0	PRO	-	expression tag	UNP Q8LIL0
B	46	ILE	PHE	SEE REMARK 999	UNP Q8LIL0
C	-13	MET	-	expression tag	UNP Q8LIL0
C	-12	GLY	-	expression tag	UNP Q8LIL0
C	-11	SER	-	expression tag	UNP Q8LIL0
C	-10	SER	-	expression tag	UNP Q8LIL0
C	-9	HIS	-	expression tag	UNP Q8LIL0
C	-8	HIS	-	expression tag	UNP Q8LIL0
C	-7	HIS	-	expression tag	UNP Q8LIL0
C	-6	HIS	-	expression tag	UNP Q8LIL0
C	-5	HIS	-	expression tag	UNP Q8LIL0
C	-4	HIS	-	expression tag	UNP Q8LIL0
C	-3	SER	-	expression tag	UNP Q8LIL0
C	-2	GLN	-	expression tag	UNP Q8LIL0
C	-1	ASP	-	expression tag	UNP Q8LIL0
C	0	PRO	-	expression tag	UNP Q8LIL0
C	46	ILE	PHE	SEE REMARK 999	UNP Q8LIL0
D	-13	MET	-	expression tag	UNP Q8LIL0
D	-12	GLY	-	expression tag	UNP Q8LIL0
D	-11	SER	-	expression tag	UNP Q8LIL0
D	-10	SER	-	expression tag	UNP Q8LIL0
D	-9	HIS	-	expression tag	UNP Q8LIL0
D	-8	HIS	-	expression tag	UNP Q8LIL0
D	-7	HIS	-	expression tag	UNP Q8LIL0
D	-6	HIS	-	expression tag	UNP Q8LIL0
D	-5	HIS	-	expression tag	UNP Q8LIL0
D	-4	HIS	-	expression tag	UNP Q8LIL0
D	-3	SER	-	expression tag	UNP Q8LIL0
D	-2	GLN	-	expression tag	UNP Q8LIL0
D	-1	ASP	-	expression tag	UNP Q8LIL0
D	0	PRO	-	expression tag	UNP Q8LIL0
D	46	ILE	PHE	SEE REMARK 999	UNP Q8LIL0

- Molecule 2 is water.

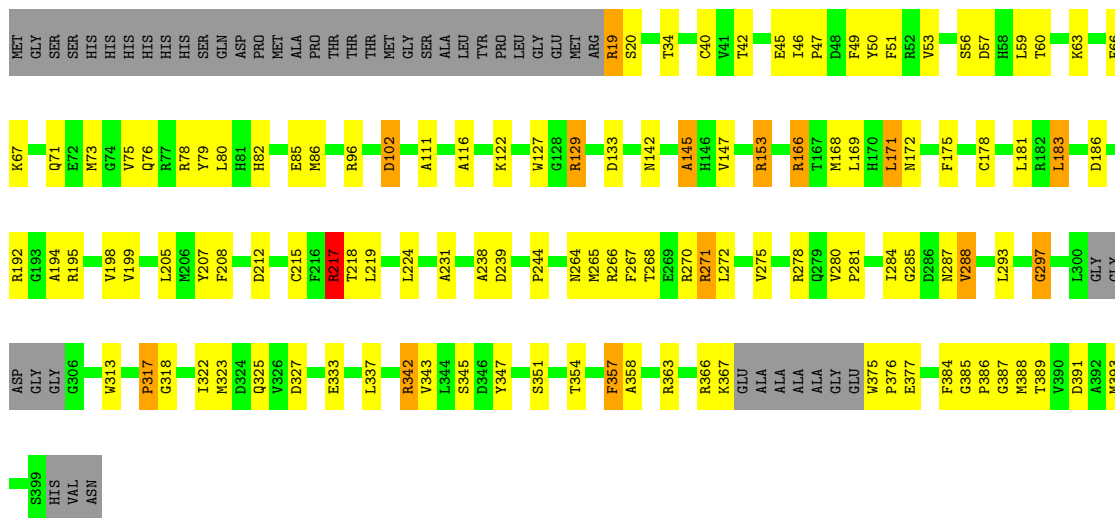
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	28	Total 28	O 28	0	0
2	B	45	Total 45	O 45	0	0
2	C	31	Total 31	O 31	0	0
2	D	21	Total 21	O 21	0	0

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.

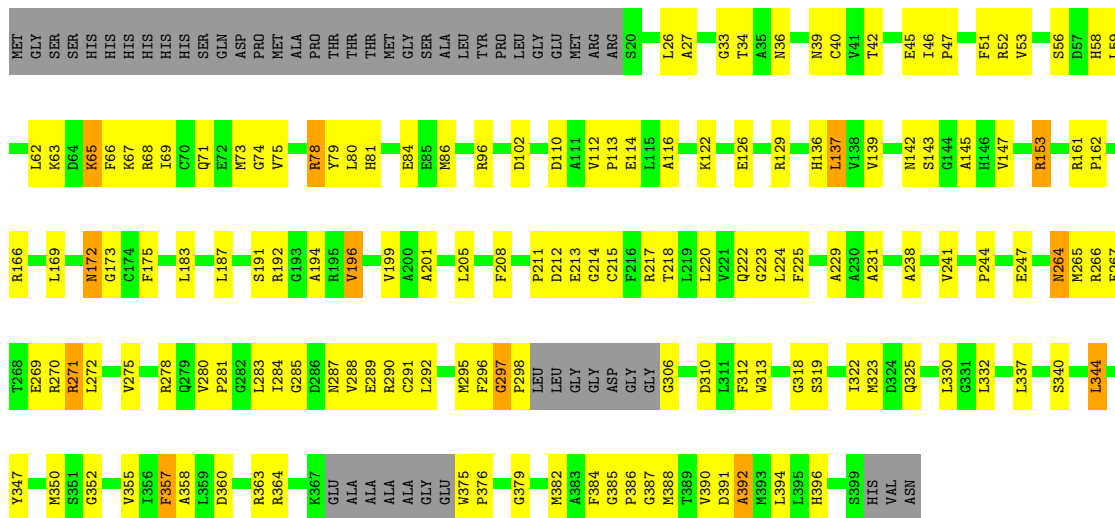
• Molecule 1: Os07g0271500 protein

Chain A: 



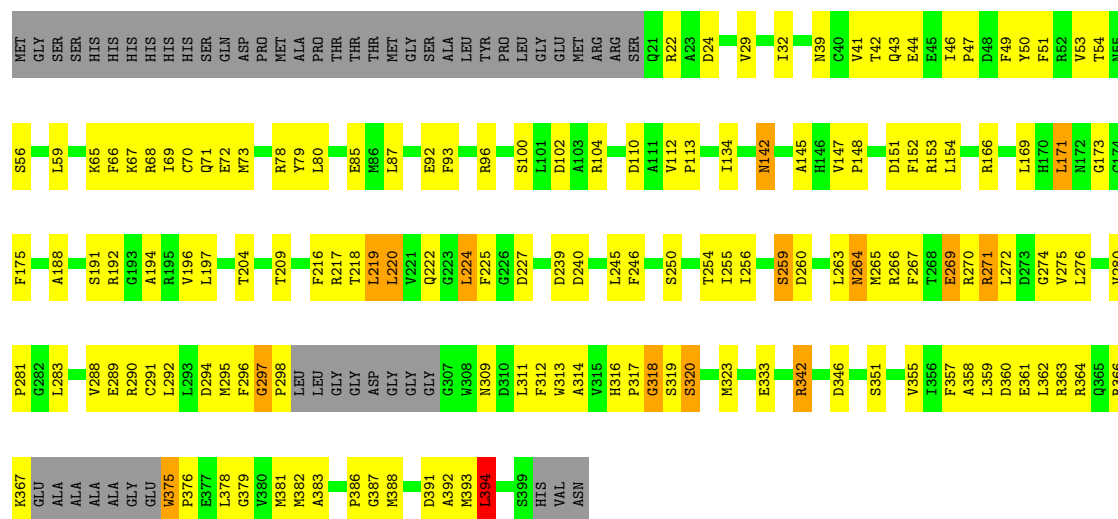
• Molecule 1: Os07g0271500 protein

Chain B: 



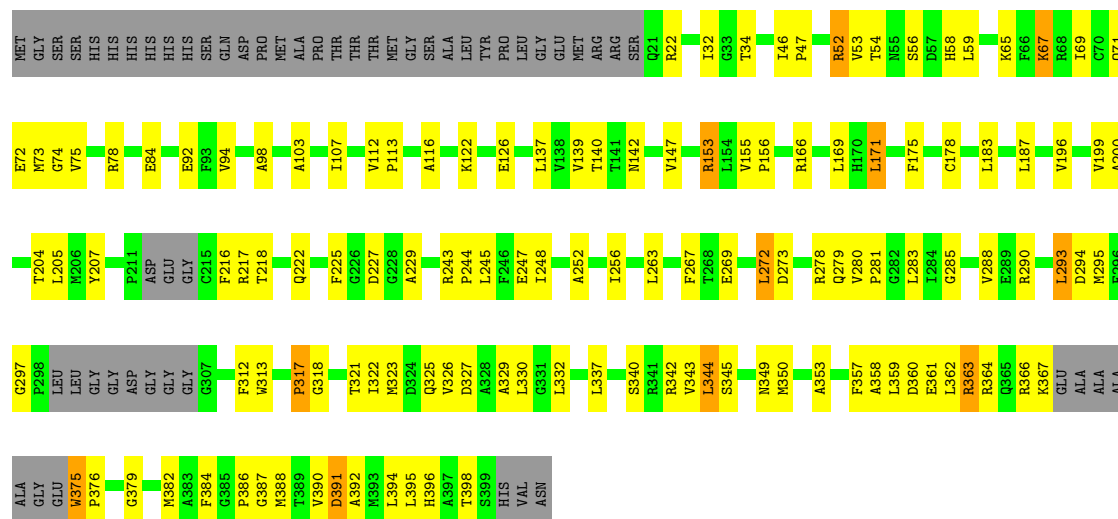
• Molecule 1: Os07g0271500 protein

Chain C:  52% 32% 12%



• Molecule 1: Os07g0271500 protein

Chain D:  56% 29% 13%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	72.70Å 97.20Å 126.20Å 90.00° 103.70° 90.00°	Depositor
Resolution (Å)	25.96 – 2.50 25.96 – 2.50	Depositor EDS
% Data completeness (in resolution range)	92.7 (25.96-2.50) 92.7 (25.96-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.03 (at 2.50Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.221 , 0.272 0.222 , 0.274	Depositor DCC
R_{free} test set	2771 reflections (5.05%)	wwPDB-VP
Wilson B-factor (Å ²)	29.7	Xtriage
Anisotropy	0.078	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 19.7	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	11223	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.88% 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	0.47	0/2858	1.02	13/3877 (0.3%)
1	B	0.52	0/2831	0.99	14/3841 (0.4%)
1	C	0.45	0/2821	0.93	14/3828 (0.4%)
1	D	0.41	0/2799	0.97	9/3797 (0.2%)
All	All	0.46	0/11309	0.98	50/15343 (0.3%)

There are no bond length outliers.

All (50) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	256	ILE	N-CA-C	-9.33	100.61	108.63
1	C	318	GLY	N-CA-C	-7.30	103.80	112.64
1	B	196	VAL	N-CA-C	7.21	118.25	107.80
1	A	129	ARG	CA-C-N	7.03	126.79	119.76
1	A	129	ARG	C-N-CA	7.03	126.79	119.76
1	A	111	ALA	N-CA-C	6.77	118.74	111.36
1	B	166	ARG	N-CA-C	6.77	120.86	109.76
1	C	240	ASP	N-CA-C	6.63	119.85	111.69
1	A	313	TRP	N-CA-C	6.61	119.93	108.76
1	A	354	THR	N-CA-C	6.37	118.76	111.11
1	D	272	LEU	N-CA-C	-6.37	99.63	109.76
1	C	313	TRP	N-CA-C	6.32	119.44	108.76
1	B	208	PHE	N-CA-C	6.32	119.38	109.96
1	C	220	LEU	N-CA-C	-6.25	104.00	111.69
1	B	357	PHE	N-CA-C	-6.16	104.57	111.28
1	B	129	ARG	CA-C-N	6.02	125.97	119.76
1	B	129	ARG	C-N-CA	6.02	125.97	119.76
1	B	313	TRP	N-CA-C	6.01	119.28	109.24
1	C	267	PHE	N-CA-C	-6.00	100.73	110.20
1	B	81	HIS	N-CA-C	-5.95	104.88	111.36
1	D	122	LYS	N-CA-C	-5.95	104.88	111.36
1	C	196	VAL	N-CA-C	5.92	116.64	107.99
1	C	357	PHE	N-CA-C	-5.92	104.95	111.82

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	166	ARG	N-CA-C	5.83	119.56	110.17
1	A	50	TYR	N-CA-C	5.82	117.30	111.07
1	C	142	ASN	N-CA-C	-5.71	106.15	113.23
1	C	394	LEU	N-CA-C	-5.69	99.92	108.96
1	C	272	LEU	N-CA-C	-5.66	100.07	109.46
1	A	239	ASP	N-CA-C	-5.58	103.34	110.53
1	A	208	PHE	N-CA-C	5.52	118.58	110.30
1	D	98	ALA	N-CA-C	5.52	116.61	109.65
1	C	224	LEU	N-CA-C	5.47	120.82	114.04
1	A	287	ASN	N-CA-C	5.45	121.34	114.31
1	B	392	ALA	N-CA-C	5.45	117.08	108.96
1	D	273	ASP	N-CA-C	5.43	116.88	109.18
1	D	394	LEU	N-CA-C	-5.38	100.41	108.96
1	A	351	SER	CB-CA-C	-5.37	110.37	116.54
1	A	357	PHE	N-CA-C	-5.35	105.44	111.28
1	A	217	ARG	N-CA-C	-5.31	105.49	111.28
1	B	27	ALA	N-CA-C	-5.28	102.46	110.28
1	B	352	GLY	N-CA-C	5.26	120.27	113.27
1	B	142	ASN	N-CA-C	-5.22	106.75	113.23
1	D	74	GLY	N-CA-C	-5.20	107.86	115.72
1	A	166	ARG	N-CA-C	5.20	118.59	110.32
1	C	351	SER	CB-CA-C	-5.17	110.60	116.54
1	D	196	VAL	N-CA-C	5.12	115.47	107.99
1	C	225	PHE	N-CA-C	5.12	118.56	109.96
1	B	36	ASN	CA-C-N	5.06	123.35	119.66
1	B	36	ASN	C-N-CA	5.06	123.35	119.66
1	C	346	ASP	N-CA-C	5.03	119.47	113.23

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	2805	0	2794	93	0
1	B	2778	0	2759	122	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	C	2768	0	2751	139	0
1	D	2747	0	2737	109	0
2	A	28	0	0	2	0
2	B	45	0	0	3	0
2	C	31	0	0	4	0
2	D	21	0	0	1	0
All	All	11223	0	11041	444	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 20.

All (444) 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:B:75:VAL:HG11	1:B:78:ARG:HD2	1.43	0.99
1:C:259:SER:HB3	1:C:388:MET:H	1.30	0.96
1:C:361:GLU:HA	1:C:364:ARG:HD2	1.49	0.94
1:D:290:ARG:HD2	1:D:293:LEU:HD11	1.51	0.92
1:D:140:THR:HG22	1:D:169:LEU:HB2	1.50	0.90
1:B:318:GLY:HA3	1:B:322:ILE:HD12	1.54	0.87
1:C:29:VAL:HG23	1:C:245:LEU:HB2	1.56	0.86
1:B:46:ILE:HG22	1:B:47:PRO:HD3	1.58	0.85
1:A:183:LEU:HD11	1:B:183:LEU:HD22	1.58	0.84
1:A:46:ILE:HG13	1:A:78:ARG:HH21	1.44	0.82
1:D:155:VAL:HB	1:D:156:PRO:HD3	1.60	0.82
1:B:46:ILE:HD13	1:B:78:ARG:HH22	1.45	0.81
1:B:284:ILE:CD1	1:B:322:ILE:HG23	2.10	0.80
1:B:75:VAL:CG1	1:B:78:ARG:HD2	2.11	0.80
1:A:217:ARG:HB3	1:A:217:ARG:HH21	1.48	0.78
1:B:284:ILE:HD13	1:B:322:ILE:HG23	1.67	0.77
1:D:46:ILE:HG22	1:D:47:PRO:HD3	1.67	0.77
1:A:147:VAL:HG12	1:B:386:PRO:HB2	1.67	0.77
1:B:73:MET:SD	1:B:224:LEU:HD13	2.24	0.76
1:C:220:LEU:HD23	1:C:276:LEU:HD13	1.68	0.76
1:D:263:LEU:HD12	1:D:388:MET:HE3	1.68	0.75
1:C:192:ARG:HG3	1:C:192:ARG:HH11	1.51	0.75
1:B:270:ARG:HB3	1:B:271:ARG:NH1	2.02	0.74
1:B:363:ARG:HD2	1:B:364:ARG:HD2	1.70	0.73
1:C:259:SER:OG	1:C:283:LEU:HD13	1.89	0.73
1:B:360:ASP:O	1:B:364:ARG:HD3	1.87	0.73
1:A:76:GLN:HG2	1:A:345:SER:O	1.89	0.72

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:323:MET:HE1	1:D:340:SER:HB2	1.70	0.72
1:C:323:MET:HA	1:C:323:MET:HE2	1.73	0.71
1:A:323:MET:HE2	1:A:323:MET:HA	1.73	0.70
1:A:194:ALA:O	1:A:195:ARG:HD3	1.91	0.70
1:C:147:VAL:HG12	1:D:386:PRO:HB2	1.73	0.70
1:D:34:THR:HB	1:D:357:PHE:CZ	2.27	0.69
1:D:342:ARG:O	1:D:342:ARG:HD3	1.92	0.69
1:B:75:VAL:HG11	1:B:78:ARG:CD	2.21	0.69
1:C:169:LEU:HD23	1:D:169:LEU:HD23	1.73	0.69
1:C:69:ILE:HD11	1:C:219:LEU:HD23	1.75	0.68
1:A:217:ARG:HB3	1:A:217:ARG:NH2	2.06	0.68
1:B:375:TRP:N	1:B:376:PRO:HD2	2.08	0.68
1:A:207:TYR:CE1	1:A:272:LEU:HD23	2.29	0.68
1:B:153:ARG:HD2	2:B:434:HOH:O	1.94	0.67
1:C:39:ASN:HB3	1:C:80:LEU:O	1.94	0.67
1:A:217:ARG:HH21	1:A:217:ARG:CB	2.08	0.66
1:B:323:MET:HE1	1:B:340:SER:HB2	1.77	0.66
1:B:284:ILE:HG13	1:B:285:GLY:N	2.09	0.66
1:C:65:LYS:HG3	1:C:68:ARG:NH1	2.09	0.66
1:C:46:ILE:HD12	1:C:78:ARG:HH21	1.60	0.66
1:C:309:ASN:OD1	1:C:333:GLU:HG3	1.95	0.66
1:D:218:THR:O	1:D:222:GLN:HG2	1.96	0.66
1:B:42:THR:OG1	1:B:45:GLU:HG2	1.95	0.66
1:C:316:HIS:HB3	1:C:383:ALA:HA	1.77	0.66
1:A:318:GLY:HA3	1:A:322:ILE:HD12	1.79	0.65
1:C:375:TRP:HA	1:C:375:TRP:CE3	2.30	0.65
1:D:245:LEU:HD13	1:D:363:ARG:NH1	2.12	0.65
1:B:78:ARG:NH1	1:B:223:GLY:HA2	2.12	0.65
1:D:323:MET:HE3	1:D:337:LEU:HB3	1.79	0.65
1:B:65:LYS:O	1:B:69:ILE:HG13	1.97	0.64
1:B:46:ILE:HD13	1:B:78:ARG:NH2	2.13	0.64
1:C:112:VAL:HB	1:C:113:PRO:HD3	1.80	0.64
1:C:265:MET:HE1	2:C:411:HOH:O	1.97	0.64
1:A:42:THR:OG1	1:A:45:GLU:HG3	1.98	0.64
1:C:46:ILE:HB	1:C:47:PRO:HD3	1.80	0.64
1:B:280:VAL:HB	1:B:281:PRO:HD3	1.79	0.63
1:D:46:ILE:CG2	1:D:47:PRO:HD3	2.29	0.63
1:D:267:PHE:HD2	1:D:272:LEU:HD13	1.63	0.63
1:D:360:ASP:O	1:D:364:ARG:HG3	1.99	0.63
1:D:46:ILE:HG12	1:D:78:ARG:HH22	1.64	0.63
1:C:204:THR:HG23	1:C:227:ASP:OD1	1.99	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:375:TRP:N	1:D:376:PRO:HD2	2.15	0.62
1:B:173:GLY:HA2	1:B:265:MET:HE1	1.82	0.62
1:C:291:CYS:O	1:C:295:MET:HG3	1.98	0.62
1:B:285:GLY:HA3	1:B:325:GLN:OE1	2.00	0.62
1:D:293:LEU:HD12	1:D:294:ASP:OD1	2.00	0.62
1:C:24:ASP:CG	1:C:192:ARG:HH22	2.08	0.61
1:B:46:ILE:CG2	1:B:47:PRO:HD3	2.28	0.61
1:B:126:GLU:CD	1:B:363:ARG:HH22	2.09	0.61
1:C:69:ILE:CD1	1:C:219:LEU:HD23	2.31	0.61
1:A:19:ARG:HA	1:B:26:LEU:HD21	1.82	0.61
1:D:183:LEU:O	1:D:187:LEU:HD13	2.00	0.61
1:A:46:ILE:HG13	1:A:78:ARG:NH2	2.15	0.61
1:C:29:VAL:HG22	1:C:246:PHE:H	1.66	0.60
1:B:47:PRO:HA	1:B:66:PHE:HE2	1.67	0.60
1:D:280:VAL:HB	1:D:281:PRO:HD3	1.83	0.60
1:C:151:ASP:OD2	1:C:166:ARG:HD2	2.02	0.60
1:B:52:ARG:HH22	1:B:84:GLU:CD	2.10	0.60
1:A:280:VAL:HB	1:A:281:PRO:HD3	1.83	0.60
1:A:284:ILE:O	1:A:288:VAL:HG22	2.02	0.59
1:B:288:VAL:HG13	1:B:289:GLU:N	2.17	0.59
1:C:192:ARG:HG3	1:C:192:ARG:NH1	2.18	0.59
1:D:384:PHE:CD1	1:D:390:VAL:HG22	2.37	0.59
1:D:387:GLY:N	1:D:388:MET:HA	2.17	0.59
1:B:112:VAL:HB	1:B:113:PRO:HD3	1.84	0.59
1:B:296:PHE:O	1:B:297:GLY:C	2.46	0.59
1:C:314:ALA:O	1:C:381:MET:HA	2.03	0.59
1:C:375:TRP:HA	1:C:375:TRP:HE3	1.66	0.59
1:C:22:ARG:HD3	1:D:22:ARG:HD3	1.83	0.58
1:C:142:ASN:OD1	1:C:171:LEU:HD12	2.03	0.58
1:A:153:ARG:HH11	1:A:153:ARG:CG	2.16	0.58
1:B:323:MET:CE	1:B:337:LEU:HB3	2.32	0.58
1:D:75:VAL:HG22	1:D:344:LEU:CD2	2.33	0.58
1:A:78:ARG:HG2	1:A:347:TYR:O	2.02	0.58
1:D:363:ARG:O	1:D:367:LYS:HG2	2.03	0.58
1:A:386:PRO:HB2	1:B:147:VAL:HG12	1.84	0.58
1:C:216:PHE:O	1:C:219:LEU:HD22	2.03	0.58
1:D:330:LEU:HD12	1:D:332:LEU:HD11	1.84	0.58
1:C:29:VAL:CG2	1:C:245:LEU:HB2	2.33	0.57
1:D:140:THR:CG2	1:D:169:LEU:HD12	2.34	0.57
1:A:102:ASP:HB2	1:B:269:GLU:OE2	2.05	0.57
1:A:363:ARG:O	1:A:367:LYS:HG2	2.04	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:278:ARG:HG2	1:D:278:ARG:HH11	1.69	0.57
1:D:140:THR:HG22	1:D:169:LEU:HD12	1.87	0.57
1:B:56:SER:HB2	1:B:59:LEU:HD12	1.87	0.57
1:C:250:SER:O	1:C:393:MET:HA	2.04	0.57
1:D:243:ARG:HG3	1:D:243:ARG:O	2.04	0.57
1:C:312:PHE:CE1	1:C:379:GLY:HA3	2.40	0.57
1:C:153:ARG:HH11	1:C:153:ARG:CG	2.18	0.57
1:B:270:ARG:HB3	1:B:271:ARG:HH12	1.67	0.56
1:C:56:SER:HB3	1:C:59:LEU:HD23	1.87	0.56
1:B:137:LEU:HD22	1:B:139:VAL:HG23	1.87	0.56
1:C:49:PHE:O	1:C:53:VAL:HG13	2.05	0.56
1:C:65:LYS:HE3	1:C:68:ARG:HH12	1.70	0.56
1:C:217:ARG:NH2	1:C:276:LEU:H	2.04	0.56
1:D:267:PHE:CD2	1:D:272:LEU:HD13	2.41	0.56
1:D:290:ARG:HH21	1:D:290:ARG:HG3	1.69	0.56
1:B:394:LEU:C	1:B:394:LEU:HD23	2.30	0.56
1:B:78:ARG:HH11	1:B:223:GLY:HA2	1.69	0.56
1:D:73:MET:HA	1:D:73:MET:HE2	1.88	0.56
1:A:82:HIS:HB3	1:A:86:MET:HE2	1.87	0.56
1:B:47:PRO:HA	1:B:66:PHE:CE2	2.41	0.56
1:A:169:LEU:HD23	1:B:169:LEU:HD22	1.88	0.56
1:C:153:ARG:HH11	1:C:153:ARG:HG3	1.70	0.56
1:C:173:GLY:HA2	1:C:265:MET:CE	2.36	0.55
1:C:288:VAL:HG13	1:C:289:GLU:N	2.22	0.55
1:A:46:ILE:HB	1:A:47:PRO:HD3	1.88	0.55
1:C:148:PRO:HB2	1:C:152:PHE:CB	2.36	0.55
1:D:103:ALA:O	1:D:107:ILE:HG13	2.07	0.55
1:D:285:GLY:HA3	1:D:325:GLN:HE21	1.71	0.55
1:D:313:TRP:O	1:D:362:LEU:HD22	2.07	0.55
1:C:79:TYR:O	1:C:227:ASP:HB2	2.07	0.55
1:D:67:LYS:O	1:D:71:GLN:HG2	2.07	0.55
1:B:205:LEU:C	1:B:205:LEU:HD12	2.32	0.54
1:D:75:VAL:HA	1:D:345:SER:HA	1.88	0.54
1:B:137:LEU:CD2	1:B:139:VAL:HG23	2.38	0.54
1:C:264:ASN:O	1:C:274:GLY:HA3	2.07	0.54
1:D:244:PRO:HB2	1:D:398:THR:HG21	1.89	0.54
1:D:112:VAL:HB	1:D:113:PRO:HD3	1.90	0.54
1:D:323:MET:CE	1:D:340:SER:HB2	2.37	0.54
1:B:288:VAL:HG13	1:B:289:GLU:H	1.73	0.54
1:C:366:ARG:HH12	1:C:376:PRO:HD2	1.73	0.54
1:A:278:ARG:HG2	1:A:278:ARG:HH11	1.73	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:168:MET:HB3	1:B:172:ASN:HD21	1.73	0.54
1:B:247:GLU:HB2	1:B:396:HIS:HB3	1.89	0.54
1:C:270:ARG:O	1:C:271:ARG:HB3	2.08	0.54
1:C:355:VAL:O	1:C:358:ALA:HB3	2.08	0.54
1:C:259:SER:OG	1:C:388:MET:HG3	2.08	0.54
1:A:56:SER:HB3	1:A:59:LEU:HD12	1.90	0.53
1:C:68:ARG:HG3	1:C:69:ILE:N	2.23	0.53
1:C:312:PHE:CD1	1:C:379:GLY:HA3	2.42	0.53
1:D:52:ARG:HD2	1:D:52:ARG:C	2.33	0.53
1:A:387:GLY:N	1:A:388:MET:HA	2.22	0.53
1:A:342:ARG:HH21	1:A:342:ARG:HG3	1.74	0.53
1:A:198:VAL:O	1:A:231:ALA:HA	2.09	0.53
1:D:285:GLY:O	1:D:329:ALA:HB2	2.08	0.53
1:A:67:LYS:O	1:A:71:GLN:HG3	2.09	0.53
1:D:343:VAL:HG21	1:D:358:ALA:HA	1.91	0.53
1:A:47:PRO:HG2	1:A:67:LYS:HD2	1.91	0.52
1:B:52:ARG:NH2	1:B:84:GLU:CD	2.67	0.52
1:D:216:PHE:CD2	1:D:278:ARG:NH2	2.77	0.52
1:D:247:GLU:HB2	1:D:396:HIS:HB3	1.91	0.52
1:C:256:ILE:O	1:C:259:SER:HB2	2.08	0.52
1:C:288:VAL:HG13	1:C:289:GLU:H	1.73	0.52
1:B:291:CYS:O	1:B:295:MET:HG3	2.08	0.52
1:C:29:VAL:HG23	1:C:29:VAL:O	2.10	0.52
1:A:366:ARG:O	1:A:366:ARG:HG2	2.10	0.52
1:C:382:MET:HB2	1:C:392:ALA:HB2	1.92	0.52
1:D:361:GLU:HB2	1:D:364:ARG:HH21	1.75	0.52
1:B:264:ASN:OD1	1:B:275:VAL:HB	2.10	0.52
1:D:318:GLY:HA2	1:D:349:ASN:ND2	2.25	0.52
1:C:173:GLY:HA2	1:C:265:MET:HE1	1.92	0.52
1:D:53:VAL:HG21	1:D:84:GLU:HA	1.92	0.52
1:D:56:SER:HB3	1:D:59:LEU:CD1	2.40	0.52
1:C:317:PRO:HA	1:C:323:MET:HE3	1.92	0.51
1:A:153:ARG:HH11	1:A:153:ARG:HG3	1.74	0.51
1:D:312:PHE:CD1	1:D:379:GLY:HA3	2.46	0.51
1:A:19:ARG:NE	1:A:20:SER:O	2.41	0.51
1:A:34:THR:O	1:A:122:LYS:HD2	2.10	0.51
1:A:270:ARG:O	1:A:271:ARG:HB3	2.11	0.51
1:B:278:ARG:HG2	1:B:278:ARG:HH11	1.76	0.51
1:C:217:ARG:HH21	1:C:217:ARG:HG3	1.74	0.51
1:B:323:MET:HE3	1:B:337:LEU:HB3	1.92	0.51
1:A:42:THR:HG23	1:A:45:GLU:OE1	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:327:ASP:OD2	1:A:337:LEU:HD12	2.11	0.51
1:D:321:THR:O	1:D:325:GLN:HG2	2.11	0.51
1:D:137:LEU:HD13	1:D:137:LEU:C	2.36	0.51
1:A:82:HIS:CB	1:A:86:MET:HE2	2.41	0.51
1:C:342:ARG:CG	1:C:342:ARG:HH21	2.23	0.51
1:D:278:ARG:HG2	1:D:278:ARG:NH1	2.26	0.51
1:B:40:CYS:HB2	1:B:79:TYR:CE1	2.46	0.51
1:B:86:MET:HE1	1:B:205:LEU:HD22	1.93	0.51
1:B:284:ILE:HD11	1:B:322:ILE:HG23	1.91	0.51
1:B:363:ARG:CD	1:B:364:ARG:HD2	2.39	0.51
1:C:265:MET:O	1:C:266:ARG:HG3	2.11	0.51
1:C:382:MET:CB	1:C:392:ALA:HB2	2.40	0.51
1:C:296:PHE:O	1:C:297:GLY:C	2.53	0.51
1:D:116:ALA:HB1	1:D:199:VAL:HG11	1.92	0.51
1:B:212:ASP:O	1:B:214:GLY:N	2.44	0.51
1:C:148:PRO:HB2	1:C:152:PHE:HB3	1.92	0.51
1:C:224:LEU:HD21	1:C:318:GLY:O	2.11	0.51
1:C:259:SER:HG	1:C:283:LEU:HD13	1.76	0.51
1:B:175:PHE:HD2	1:B:385:GLY:HA3	1.75	0.50
1:D:92:GLU:H	1:D:92:GLU:CD	2.19	0.50
1:B:217:ARG:HG2	1:B:217:ARG:NH2	2.25	0.50
1:D:322:ILE:N	1:D:322:ILE:HD12	2.26	0.50
1:B:173:GLY:CA	1:B:265:MET:HE1	2.40	0.50
1:D:140:THR:HG22	1:D:169:LEU:CB	2.32	0.50
1:A:169:LEU:HD23	1:B:169:LEU:CD2	2.42	0.50
1:A:215:CYS:SG	1:A:217:ARG:CZ	2.99	0.50
1:B:110:ASP:O	1:B:114:GLU:HG2	2.11	0.50
1:D:363:ARG:HG2	1:D:363:ARG:HH11	1.76	0.50
1:B:267:PHE:CD1	1:B:272:LEU:HD13	2.47	0.50
1:B:217:ARG:HG2	1:B:217:ARG:HH21	1.77	0.50
1:C:260:ASP:HA	2:C:418:HOH:O	2.11	0.50
1:D:56:SER:HB3	1:D:59:LEU:HD12	1.94	0.50
1:B:73:MET:HE2	1:B:73:MET:HA	1.93	0.50
1:B:53:VAL:HG12	1:B:84:GLU:HG3	1.93	0.49
1:D:67:LYS:HE3	1:D:71:GLN:NE2	2.27	0.49
1:A:215:CYS:HB3	2:A:414:HOH:O	2.12	0.49
1:C:290:ARG:NE	1:C:294:ASP:OD2	2.45	0.49
1:A:34:THR:HB	1:A:357:PHE:CZ	2.47	0.49
1:B:387:GLY:N	1:B:388:MET:HA	2.27	0.49
1:C:216:PHE:HA	1:C:219:LEU:CD1	2.43	0.49
1:B:238:ALA:CB	1:B:244:PRO:HG3	2.42	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:317:PRO:O	1:D:318:GLY:C	2.56	0.48
1:B:34:THR:O	1:B:122:LYS:HD2	2.13	0.48
1:B:323:MET:HE1	1:B:337:LEU:HB3	1.95	0.48
1:C:387:GLY:N	1:C:388:MET:HA	2.28	0.48
1:D:200:ALA:O	1:D:229:ALA:HA	2.13	0.48
1:B:220:LEU:O	1:B:224:LEU:HB2	2.13	0.48
1:A:80:LEU:HD22	1:A:82:HIS:CE1	2.48	0.48
1:C:280:VAL:HG13	1:C:388:MET:HE1	1.94	0.48
1:D:361:GLU:OE1	1:D:364:ARG:NH2	2.47	0.48
1:D:323:MET:HE1	1:D:340:SER:CB	2.41	0.48
1:D:332:LEU:HD13	1:D:337:LEU:CD1	2.43	0.48
1:A:142:ASN:OD1	1:A:171:LEU:HD12	2.14	0.48
1:D:327:ASP:OD1	1:D:337:LEU:HD22	2.14	0.48
1:B:312:PHE:CD1	1:B:379:GLY:HA3	2.49	0.48
1:B:323:MET:CE	1:B:340:SER:HB2	2.44	0.48
1:A:40:CYS:HB2	1:A:79:TYR:CE1	2.49	0.48
1:B:296:PHE:O	1:B:298:PRO:N	2.46	0.48
1:B:312:PHE:CE1	1:B:379:GLY:HA3	2.49	0.47
1:D:353:ALA:O	1:D:357:PHE:HD1	1.96	0.47
1:D:375:TRP:CE3	1:D:375:TRP:HA	2.49	0.47
1:B:65:LYS:HD3	1:B:68:ARG:NH2	2.30	0.47
1:C:53:VAL:HB	1:C:87:LEU:HD13	1.96	0.47
1:D:361:GLU:HA	1:D:364:ARG:HE	1.79	0.47
1:A:266:ARG:HH21	1:A:266:ARG:HG2	1.79	0.47
1:C:78:ARG:HD3	1:C:78:ARG:N	2.29	0.47
1:C:42:THR:HB	1:C:44:GLU:OE2	2.14	0.47
1:C:153:ARG:CG	1:C:153:ARG:NH1	2.77	0.47
1:B:283:LEU:O	1:B:287:ASN:ND2	2.42	0.47
1:A:47:PRO:HG2	1:A:67:LYS:CD	2.44	0.47
1:C:151:ASP:CG	1:C:152:PHE:N	2.72	0.47
1:A:266:ARG:O	1:A:272:LEU:HD12	2.15	0.47
1:B:191:SER:HB2	1:B:194:ALA:HB2	1.96	0.47
1:B:375:TRP:N	1:B:376:PRO:CD	2.76	0.47
1:C:102:ASP:OD2	1:D:269:GLU:HG3	2.14	0.47
1:C:147:VAL:HA	1:C:148:PRO:C	2.39	0.47
1:D:322:ILE:O	1:D:326:VAL:HG23	2.14	0.47
1:D:332:LEU:HD13	1:D:337:LEU:HD11	1.97	0.47
1:A:205:LEU:HD12	1:A:205:LEU:C	2.40	0.47
1:B:62:LEU:HD21	1:B:211:PRO:HB2	1.96	0.47
1:A:278:ARG:HG2	1:A:278:ARG:NH1	2.29	0.47
1:B:330:LEU:HB2	1:B:332:LEU:HG	1.97	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:263:LEU:HD23	1:C:275:VAL:O	2.15	0.47
1:C:366:ARG:NH1	1:C:376:PRO:HD2	2.30	0.47
1:A:264:ASN:OD1	1:A:275:VAL:HB	2.15	0.46
1:A:75:VAL:HA	1:A:345:SER:HA	1.96	0.46
1:A:317:PRO:O	1:A:318:GLY:C	2.58	0.46
1:C:191:SER:HB2	1:C:194:ALA:HB2	1.96	0.46
1:A:217:ARG:HG3	1:A:218:THR:N	2.30	0.46
1:B:201:ALA:HA	1:B:229:ALA:HA	1.97	0.46
1:C:46:ILE:HD11	1:C:78:ARG:HE	1.80	0.46
1:D:375:TRP:HA	1:D:375:TRP:HE3	1.79	0.46
1:A:153:ARG:CG	1:A:153:ARG:NH1	2.77	0.46
1:C:32:ILE:CD1	1:C:359:LEU:HD12	2.46	0.46
1:C:363:ARG:HD3	1:C:367:LYS:NZ	2.30	0.46
1:D:65:LYS:O	1:D:69:ILE:HG13	2.15	0.46
1:D:137:LEU:HD11	1:D:139:VAL:HG23	1.98	0.46
1:A:186:ASP:OD1	1:B:187:LEU:HD23	2.15	0.46
1:A:375:TRP:O	1:A:376:PRO:C	2.59	0.46
1:D:205:LEU:C	1:D:205:LEU:HD12	2.40	0.46
1:C:254:THR:OG1	1:C:255:ILE:N	2.48	0.46
1:D:290:ARG:NH2	1:D:294:ASP:OD1	2.48	0.46
1:A:42:THR:HG23	1:A:45:GLU:CD	2.41	0.46
1:A:280:VAL:HG13	1:A:388:MET:HE1	1.97	0.46
1:C:92:GLU:HG3	1:C:100:SER:HB3	1.98	0.46
1:A:51:PHE:CZ	1:A:63:LYS:HA	2.50	0.46
1:A:145:ALA:O	1:B:266:ARG:HA	2.16	0.46
1:B:218:THR:O	1:B:222:GLN:HG2	2.15	0.46
1:C:41:VAL:CG1	1:C:46:ILE:HD13	2.45	0.46
1:B:74:GLY:O	1:B:344:LEU:HD23	2.15	0.46
1:A:142:ASN:ND2	1:A:265:MET:HE1	2.31	0.45
1:C:73:MET:HE2	1:C:73:MET:HA	1.98	0.45
1:C:188:ALA:HA	1:C:194:ALA:HB3	1.99	0.45
1:B:51:PHE:CZ	1:B:63:LYS:HA	2.51	0.45
1:D:252:ALA:HB3	1:D:295:MET:HB3	1.97	0.45
1:B:39:ASN:HB3	1:B:80:LEU:O	2.17	0.45
1:C:73:MET:O	1:C:320:SER:HB2	2.17	0.45
1:C:175:PHE:C	1:C:175:PHE:CD1	2.95	0.45
1:B:56:SER:HA	1:B:58:HIS:CE1	2.51	0.45
1:B:96:ARG:NH2	1:B:215:CYS:SG	2.90	0.45
1:C:169:LEU:HD13	2:C:428:HOH:O	2.16	0.45
1:C:360:ASP:OD2	1:C:364:ARG:NH2	2.50	0.45
1:D:56:SER:HA	1:D:58:HIS:CE1	2.52	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:207:TYR:HE2	2:D:419:HOH:O	1.99	0.45
1:B:306:GLY:HA2	1:B:310:ASP:OD1	2.17	0.45
1:B:355:VAL:O	1:B:358:ALA:HB3	2.16	0.45
1:C:362:LEU:O	1:C:366:ARG:HG3	2.16	0.45
1:D:142:ASN:OD1	1:D:171:LEU:HD12	2.16	0.45
1:C:67:LYS:O	1:C:71:GLN:HG3	2.16	0.45
1:C:110:ASP:O	1:C:113:PRO:HD2	2.16	0.45
1:C:342:ARG:CG	1:C:342:ARG:NH2	2.80	0.45
1:D:288:VAL:HG23	1:D:384:PHE:HZ	1.82	0.45
1:A:96:ARG:NE	1:A:212:ASP:HB2	2.32	0.44
1:B:270:ARG:O	1:B:271:ARG:HB3	2.16	0.44
1:C:43:GLN:HA	1:C:46:ILE:HG12	1.98	0.44
1:B:33:GLY:O	1:B:231:ALA:N	2.49	0.44
1:B:78:ARG:HH11	1:B:223:GLY:C	2.25	0.44
1:B:384:PHE:HD1	1:B:390:VAL:HG22	1.82	0.44
1:A:96:ARG:CZ	1:A:212:ASP:HB2	2.47	0.44
1:C:68:ARG:HG3	1:C:69:ILE:HG23	1.98	0.44
1:B:65:LYS:HB2	1:B:65:LYS:NZ	2.33	0.44
1:B:78:ARG:HH11	1:B:223:GLY:CA	2.30	0.44
1:C:32:ILE:HD11	1:C:359:LEU:HD12	1.99	0.44
1:C:59:LEU:N	1:C:59:LEU:HD22	2.33	0.44
1:C:216:PHE:HA	1:C:219:LEU:HD13	1.99	0.44
1:C:379:GLY:O	1:C:394:LEU:HD23	2.18	0.44
1:A:116:ALA:HB1	1:A:199:VAL:HG11	1.98	0.44
1:C:292:LEU:HA	1:C:295:MET:HE3	2.00	0.44
1:D:217:ARG:HG2	1:D:217:ARG:HH21	1.82	0.44
1:D:225:PHE:HD1	1:D:349:ASN:HB3	1.83	0.44
1:D:330:LEU:HB2	1:D:332:LEU:HG	2.00	0.44
1:D:363:ARG:NH1	1:D:363:ARG:HG2	2.33	0.44
1:C:216:PHE:O	1:C:217:ARG:C	2.60	0.43
1:D:342:ARG:HG2	1:D:342:ARG:HH21	1.82	0.43
1:D:366:ARG:NH1	1:D:375:TRP:N	2.66	0.43
1:D:312:PHE:CE1	1:D:379:GLY:HA3	2.52	0.43
1:A:285:GLY:HA3	1:A:325:GLN:OE1	2.18	0.43
1:C:92:GLU:HG2	1:C:93:PHE:N	2.32	0.43
1:C:311:LEU:HA	1:C:378:LEU:O	2.18	0.43
1:B:162:PRO:HD2	2:B:438:HOH:O	2.18	0.43
1:C:100:SER:O	1:C:104:ARG:HG3	2.18	0.43
1:C:269:GLU:H	1:C:269:GLU:HG2	1.49	0.43
1:C:41:VAL:HG12	1:C:46:ILE:HD13	2.00	0.43
1:C:362:LEU:O	1:C:366:ARG:CG	2.66	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:178:CYS:HB3	1:A:391:ASP:OD2	2.18	0.43
1:B:284:ILE:CG1	1:B:285:GLY:N	2.79	0.43
1:C:317:PRO:O	1:C:318:GLY:C	2.61	0.43
1:A:147:VAL:HG12	1:B:386:PRO:CB	2.44	0.43
1:A:375:TRP:N	1:A:376:PRO:CD	2.81	0.43
1:B:290:ARG:HB3	1:B:290:ARG:NH2	2.34	0.43
1:C:134:ILE:HD13	1:C:197:LEU:HB2	2.01	0.43
1:C:22:ARG:HB2	1:D:22:ARG:NH1	2.33	0.43
1:C:72:GLU:O	1:C:73:MET:HE2	2.19	0.43
1:C:110:ASP:HB3	2:C:405:HOH:O	2.17	0.43
1:D:155:VAL:CB	1:D:156:PRO:HD3	2.39	0.43
1:D:137:LEU:CD1	1:D:139:VAL:HG23	2.49	0.43
1:A:268:THR:HA	1:B:102:ASP:OD2	2.19	0.42
1:A:270:ARG:HD2	1:A:270:ARG:N	2.34	0.42
1:A:278:ARG:O	1:A:281:PRO:HD2	2.19	0.42
1:C:154:LEU:HD23	1:C:154:LEU:HA	1.92	0.42
1:A:363:ARG:O	1:A:367:LYS:HE3	2.19	0.42
1:C:224:LEU:HD21	1:C:319:SER:HB3	2.01	0.42
1:C:246:PHE:CD2	1:C:359:LEU:HD11	2.54	0.42
1:C:360:ASP:O	1:C:364:ARG:HG3	2.19	0.42
1:C:375:TRP:N	1:C:376:PRO:CD	2.83	0.42
1:A:49:PHE:O	1:A:53:VAL:HG22	2.19	0.42
1:A:73:MET:HE2	1:A:73:MET:HA	2.00	0.42
1:D:382:MET:HB3	1:D:392:ALA:HB2	2.01	0.42
1:A:384:PHE:HA	1:A:389:THR:O	2.19	0.42
1:A:34:THR:HB	1:A:357:PHE:CE2	2.54	0.42
1:C:65:LYS:NZ	1:C:216:PHE:CE2	2.79	0.42
1:C:280:VAL:HB	1:C:281:PRO:HD3	2.01	0.42
1:B:116:ALA:HB1	1:B:199:VAL:HG11	2.01	0.42
1:B:136:HIS:HB2	1:B:196:VAL:HG22	2.01	0.42
1:B:347:TYR:HB2	1:B:350:MET:HE2	2.02	0.42
1:C:96:ARG:HH12	1:C:218:THR:HG21	1.84	0.42
1:B:290:ARG:NH2	1:B:290:ARG:CB	2.82	0.42
1:A:175:PHE:HD2	1:A:385:GLY:HA3	1.84	0.42
1:A:238:ALA:HB2	1:A:244:PRO:HG3	2.02	0.42
1:B:225:PHE:N	1:B:225:PHE:CD1	2.88	0.42
1:C:216:PHE:HB3	1:C:219:LEU:HD21	2.01	0.42
1:C:218:THR:O	1:C:222:GLN:HG2	2.20	0.42
1:C:290:ARG:HD3	1:C:291:CYS:N	2.35	0.42
1:B:394:LEU:C	1:B:394:LEU:CD2	2.93	0.41
1:C:29:VAL:HG22	1:C:246:PHE:N	2.34	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:327:ASP:CG	1:D:337:LEU:HD22	2.44	0.41
1:A:129:ARG:NH1	1:A:133:ASP:OD1	2.52	0.41
1:C:342:ARG:HH21	1:C:342:ARG:HG3	1.84	0.41
1:D:245:LEU:HD13	1:D:363:ARG:HH12	1.82	0.41
1:D:279:GLN:O	1:D:283:LEU:HD23	2.20	0.41
1:A:145:ALA:HB1	1:B:267:PHE:CD2	2.55	0.41
1:B:67:LYS:O	1:B:71:GLN:NE2	2.54	0.41
1:B:290:ARG:HB3	1:B:290:ARG:HH21	1.84	0.41
1:C:366:ARG:NH1	1:C:375:TRP:N	2.68	0.41
1:B:192:ARG:HD2	2:B:412:HOH:O	2.20	0.41
1:C:32:ILE:HG21	1:C:360:ASP:HB2	2.01	0.41
1:C:66:PHE:CZ	1:C:70:CYS:SG	3.13	0.41
1:A:171:LEU:C	1:A:172:ASN:HD22	2.29	0.41
1:A:293:LEU:O	1:A:297:GLY:N	2.52	0.41
1:B:172:ASN:HD22	1:B:172:ASN:HA	1.60	0.41
1:B:278:ARG:HG2	1:B:278:ARG:NH1	2.35	0.41
1:D:248:ILE:HA	1:D:395:LEU:HD23	2.01	0.41
1:D:350:MET:HE1	1:D:357:PHE:CD2	2.56	0.41
1:D:54:THR:HG22	1:D:94:VAL:HG13	2.01	0.41
1:D:342:ARG:HD3	1:D:342:ARG:C	2.45	0.41
1:A:224:LEU:HD12	1:A:224:LEU:O	2.21	0.41
1:C:386:PRO:HB2	1:D:147:VAL:HG12	2.01	0.41
1:A:166:ARG:HH11	1:A:166:ARG:HD2	1.77	0.41
1:A:267:PHE:CD1	1:A:272:LEU:HD13	2.55	0.41
1:B:292:LEU:HA	1:B:295:MET:HE3	2.02	0.41
1:B:319:SER:H	1:B:322:ILE:HB	1.85	0.41
1:C:51:PHE:HD2	1:C:56:SER:O	2.04	0.41
1:D:32:ILE:HG21	1:D:360:ASP:HB2	2.02	0.41
1:D:153:ARG:H	1:D:153:ARG:HG2	1.71	0.41
1:A:66:PHE:CD1	1:A:219:LEU:HD22	2.56	0.41
1:A:181:LEU:HD12	1:A:393:MET:HE1	2.02	0.41
1:B:360:ASP:OD1	1:B:364:ARG:NE	2.53	0.41
1:C:65:LYS:O	1:C:68:ARG:HG2	2.21	0.41
1:C:96:ARG:HG2	1:C:209:THR:OG1	2.20	0.41
1:C:110:ASP:C	1:C:113:PRO:HD2	2.46	0.40
1:C:263:LEU:HD23	1:C:263:LEU:HA	1.93	0.40
1:A:343:VAL:HG21	1:A:358:ALA:HA	2.03	0.40
1:C:296:PHE:O	1:C:298:PRO:N	2.54	0.40
1:D:204:THR:HG23	1:D:227:ASP:OD2	2.21	0.40
1:D:285:GLY:HA3	1:D:325:GLN:NE2	2.34	0.40
1:A:217:ARG:HG2	2:A:414:HOH:O	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:ARG:HA	1:B:162:PRO:HD3	1.95	0.40
1:B:382:MET:HB2	1:B:392:ALA:HB2	2.03	0.40
1:D:175:PHE:HB2	1:D:391:ASP:HB2	2.02	0.40
1:A:127:TRP:CE2	1:A:129:ARG:HB2	2.57	0.40
1:B:34:THR:HB	1:B:357:PHE:CZ	2.56	0.40
1:C:50:TYR:CE1	1:C:54:THR:HG21	2.57	0.40
1:D:178:CYS:HB3	1:D:391:ASP:HB3	2.03	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	363/416 (87%)	346 (95%)	12 (3%)	5 (1%)	9	17
1	B	360/416 (86%)	340 (94%)	16 (4%)	4 (1%)	11	22
1	C	358/416 (86%)	334 (93%)	19 (5%)	5 (1%)	9	17
1	D	353/416 (85%)	329 (93%)	21 (6%)	3 (1%)	16	31
All	All	1434/1664 (86%)	1349 (94%)	68 (5%)	17 (1%)	10	20

All (17) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	213	GLU
1	B	297	GLY
1	A	145	ALA
1	A	271	ARG
1	A	297	GLY
1	B	145	ALA
1	C	145	ALA
1	C	271	ARG

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Mol	Chain	Res	Type
1	C	297	GLY
1	B	271	ARG
1	D	171	LEU
1	A	171	LEU
1	A	317	PRO
1	C	171	LEU
1	C	239	ASP
1	D	297	GLY
1	D	317	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	291/324 (90%)	278 (96%)	13 (4%)	24	49
1	B	288/324 (89%)	278 (96%)	10 (4%)	32	58
1	C	287/324 (89%)	277 (96%)	10 (4%)	32	58
1	D	285/324 (88%)	274 (96%)	11 (4%)	28	55
All	All	1151/1296 (89%)	1107 (96%)	44 (4%)	29	56

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	ARG
1	A	57	ASP
1	A	60	THR
1	A	85	GLU
1	A	102	ASP
1	A	153	ARG
1	A	183	LEU
1	A	192	ARG
1	A	217	ARG
1	A	288	VAL
1	A	333	GLU
1	A	342	ARG

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Mol	Chain	Res	Type
1	A	377	GLU
1	B	65	LYS
1	B	78	ARG
1	B	137	LEU
1	B	143	SER
1	B	153	ARG
1	B	172	ASN
1	B	241	VAL
1	B	264	ASN
1	B	344	LEU
1	B	391	ASP
1	C	85	GLU
1	C	219	LEU
1	C	259	SER
1	C	264	ASN
1	C	269	GLU
1	C	320	SER
1	C	342	ARG
1	C	375	TRP
1	C	391	ASP
1	C	394	LEU
1	D	52	ARG
1	D	67	LYS
1	D	72	GLU
1	D	126	GLU
1	D	153	ARG
1	D	293	LEU
1	D	344	LEU
1	D	359	LEU
1	D	363	ARG
1	D	375	TRP
1	D	391	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (20) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	21	GLN
1	A	55	ASN
1	A	76	GLN
1	A	172	ASN
1	A	222	GLN
1	A	279	GLN

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Mol	Chain	Res	Type
1	A	365	GLN
1	B	90	HIS
1	B	172	ASN
1	B	222	GLN
1	C	71	GLN
1	C	82	HIS
1	C	90	HIS
1	C	172	ASN
1	C	222	GLN
1	C	325	GLN
1	D	21	GLN
1	D	90	HIS
1	D	261	HIS
1	D	325	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](#)

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 [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	369/416 (88%)	-1.68	0 100 100	8, 21, 44, 71	0
1	B	366/416 (87%)	-1.67	0 100 100	9, 21, 41, 67	0
1	C	364/416 (87%)	-1.60	0 100 100	16, 30, 52, 75	0
1	D	361/416 (86%)	-1.58	0 100 100	17, 35, 58, 72	0
All	All	1460/1664 (87%)	-1.63	0 100 100	8, 27, 52, 75	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](#)

There are no ligands in this entry.

6.5 Other polymers [i](#)

There are no such residues in this entry.