



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 4, 2026 – 09:28 PM UTC

PDB ID : 1T4T / pdb_00001t4t
Title : arginase-dinor-NOHA complex
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Deposited on : 2004-04-30
Resolution : 2.20 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)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

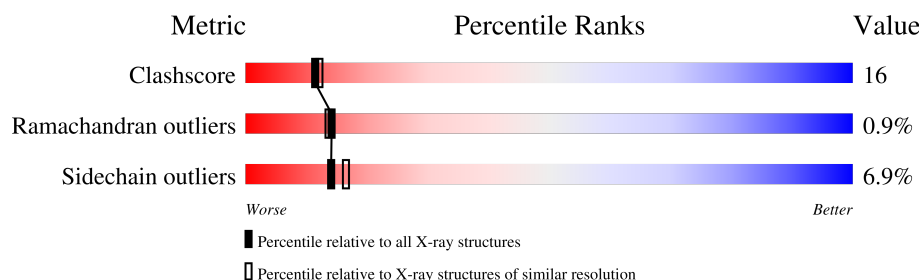
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.20 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))
Clashscore	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	314	71% 23% 6%
1	B	314	70% 25% 5%
1	C	314	69% 24% 7%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7404 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.

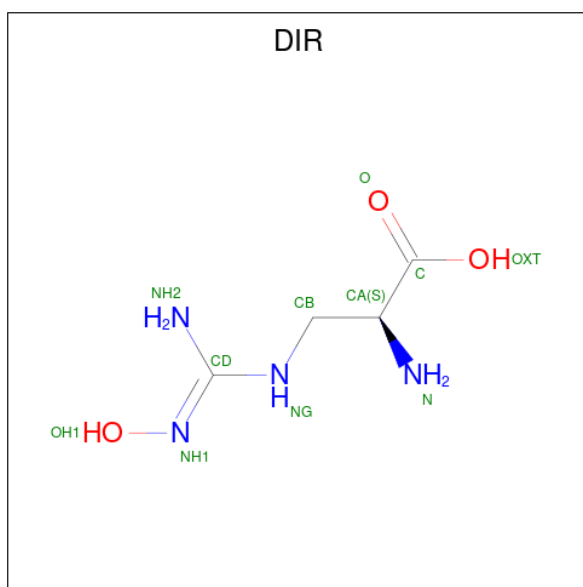
- Molecule 1 is a protein called Arginase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	314	Total	C	N	O	S	0	0	0
			2396	1528	405	456	7			
1	B	314	Total	C	N	O	S	0	0	0
			2396	1528	405	456	7			
1	C	314	Total	C	N	O	S	0	0	0
			2396	1528	405	456	7			

- Molecule 2 is MANGANESE (II) ION (CCD ID: MN) (formula: Mn).

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

- Molecule 3 is 3-[(E)-AMINO(HYDROXYIMINO)METHYL]AMINO}ALANINE (CCD ID: DIR) (formula: C₄H₁₀N₄O₃).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			11	4	4	3		
3	B	1	Total	C	N	O	0	0
			11	4	4	3		
3	C	1	Total	C	N	O	0	0
			11	4	4	3		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	44	Total	O	0	0
			44	44		
4	B	72	Total	O	0	0
			72	72		
4	C	61	Total	O	0	0
			61	61		

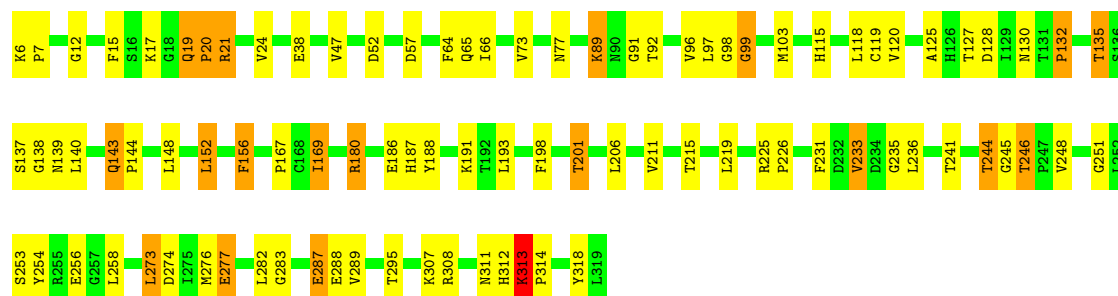
3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

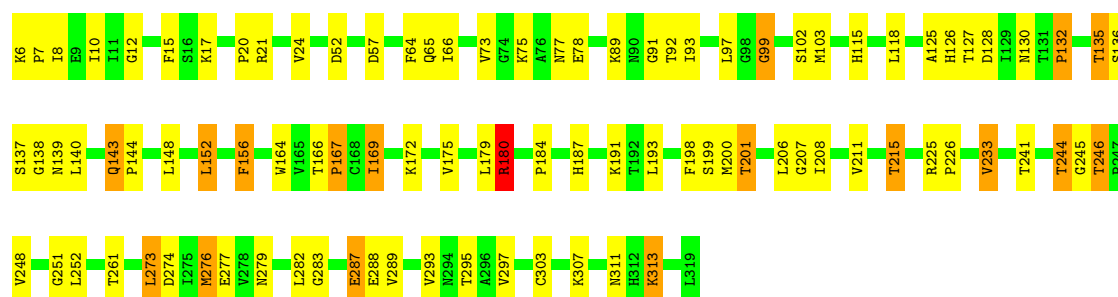
• Molecule 1: Arginase 1

Chain A: 



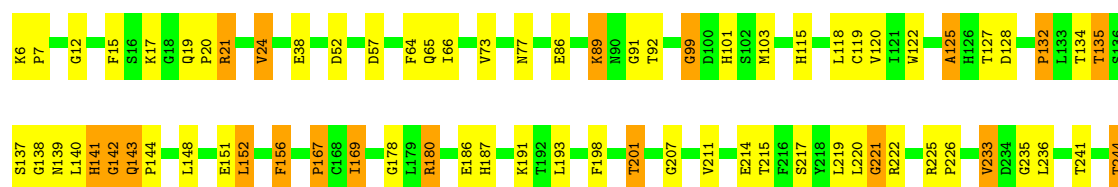
• Molecule 1: Arginase 1

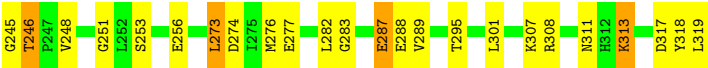
Chain B: 



• Molecule 1: Arginase 1

Chain C: 





4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 32	Depositor
Cell constants a, b, c, α , β , γ	88.30Å 88.30Å 113.10Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	28.90 – 2.20	Depositor
% Data completeness (in resolution range)	98.3 (28.90-2.20)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.258 , 0.281	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	7404	wwPDB-VP
Average B, all atoms (Å ²)	39.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: DIR, MN

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	0.51	0/2449	1.02	18/3325 (0.5%)
1	B	0.58	1/2449 (0.0%)	1.01	13/3325 (0.4%)
1	C	0.66	3/2449 (0.1%)	1.05	16/3325 (0.5%)
All	All	0.59	4/7347 (0.1%)	1.03	47/9975 (0.5%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	142	GLY	N-CA	9.81	1.59	1.45
1	C	141	HIS	C-O	-8.91	1.12	1.24
1	B	180	ARG	CZ-NH1	-5.87	1.24	1.32
1	C	135	THR	CB-CG2	-5.82	1.33	1.52

All (47) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	141	HIS	O-C-N	7.72	132.78	122.43
1	B	125	ALA	N-CA-C	-7.49	104.12	113.18
1	A	98	GLY	N-CA-C	7.34	122.94	112.81
1	A	99	GLY	N-CA-C	-7.10	96.34	113.18
1	C	99	GLY	N-CA-C	-7.06	96.45	113.18
1	B	99	GLY	N-CA-C	-7.04	96.51	113.18
1	A	125	ALA	N-CA-C	-7.00	104.71	113.18
1	B	136	SER	N-CA-C	-6.99	104.57	113.23
1	B	207	GLY	N-CA-C	-6.70	103.20	112.18
1	A	276	MET	N-CA-C	6.63	119.89	110.14
1	C	19	GLN	N-CA-C	6.62	113.96	108.07
1	C	276	MET	N-CA-C	6.58	119.81	110.14
1	A	97	LEU	CA-C-N	-6.41	116.02	122.69
1	A	97	LEU	C-N-CA	-6.41	116.02	122.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	134	THR	CB-CA-C	-6.38	103.17	111.40
1	C	207	GLY	N-CA-C	-6.31	103.72	112.18
1	B	206	LEU	N-CA-C	6.19	119.30	111.69
1	B	276	MET	N-CA-C	6.12	119.18	109.81
1	A	120	VAL	N-CA-C	6.09	116.64	108.11
1	C	134	THR	CA-C-N	-6.00	113.01	122.73
1	C	134	THR	C-N-CA	-6.00	113.01	122.73
1	C	19	GLN	CA-C-N	5.98	127.31	119.84
1	C	19	GLN	C-N-CA	5.98	127.31	119.84
1	C	125	ALA	N-CA-C	-5.91	106.03	113.18
1	A	236	LEU	N-CA-C	-5.69	102.10	110.52
1	B	180	ARG	NE-CZ-NH2	5.61	124.25	119.20
1	B	21	ARG	N-CA-C	5.37	117.46	109.25
1	C	313	LYS	CA-C-N	5.34	125.28	119.78
1	C	313	LYS	C-N-CA	5.34	125.28	119.78
1	B	313	LYS	CA-C-N	5.33	125.27	119.78
1	B	313	LYS	C-N-CA	5.33	125.27	119.78
1	A	277	GLU	N-CA-C	5.31	119.34	112.86
1	C	21	ARG	N-CA-C	5.31	117.86	109.96
1	C	217	SER	N-CA-C	-5.30	105.62	111.71
1	A	206	LEU	N-CA-C	5.26	118.16	111.69
1	A	21	ARG	N-CA-C	5.23	117.25	109.25
1	A	313	LYS	CA-C-N	5.19	125.13	119.78
1	A	313	LYS	C-N-CA	5.19	125.13	119.78
1	A	231	PHE	N-CA-C	5.18	117.07	108.20
1	B	279	ASN	CA-C-N	5.18	126.31	119.84
1	B	279	ASN	C-N-CA	5.18	126.31	119.84
1	B	97	LEU	N-CA-C	-5.16	100.75	108.96
1	A	47	VAL	N-CA-C	5.12	115.28	108.11
1	A	19	GLN	CA-C-N	5.10	124.89	119.28
1	A	19	GLN	C-N-CA	5.10	124.89	119.28
1	A	20	PRO	N-CA-C	5.08	120.50	113.65
1	C	120	VAL	N-CA-C	5.06	115.77	108.48

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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	2396	0	2420	77	0
1	B	2396	0	2420	85	0
1	C	2396	0	2420	79	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
3	A	11	0	8	1	0
3	B	11	0	8	2	0
3	C	11	0	8	2	0
4	A	44	0	0	5	0
4	B	72	0	0	8	0
4	C	61	0	0	4	0
All	All	7404	0	7284	233	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 16.

All (233) 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:132:PRO:HD2	1:B:156:PHE:HD2	1.12	1.09
1:C:132:PRO:HD2	1:C:156:PHE:HD2	1.19	1.07
1:A:132:PRO:HD2	1:A:156:PHE:HD2	1.22	1.04
1:A:7:PRO:HB2	1:A:92:THR:HG22	1.40	1.03
1:B:7:PRO:HB2	1:B:92:THR:HG22	1.41	1.02
1:C:7:PRO:HB2	1:C:92:THR:HG22	1.44	0.97
1:B:132:PRO:HD2	1:B:156:PHE:CD2	2.02	0.94
1:A:132:PRO:HD2	1:A:156:PHE:CD2	2.04	0.91
1:C:211:VAL:O	1:C:215:THR:HG23	1.71	0.91
1:C:132:PRO:HD2	1:C:156:PHE:CD2	2.07	0.89
1:C:143:GLN:N	1:C:144:PRO:HD3	1.92	0.85
1:A:211:VAL:O	1:A:215:THR:HG23	1.77	0.84
1:A:143:GLN:N	1:A:144:PRO:HD3	1.94	0.82
1:A:128:ASP:HB3	1:A:144:PRO:HD2	1.60	0.82
1:B:143:GLN:N	1:B:144:PRO:HD3	1.95	0.81
1:C:128:ASP:HB3	1:C:144:PRO:HD2	1.64	0.80
1:C:307:LYS:H	1:C:311:ASN:HD21	1.31	0.79
1:B:307:LYS:H	1:B:311:ASN:HD21	1.31	0.79
1:B:128:ASP:HB3	1:B:144:PRO:HD2	1.65	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:198:PHE:CE2	1:C:215:THR:HG22	2.18	0.78
1:B:180:ARG:NH1	1:B:251:GLY:CA	2.48	0.77
1:A:246:THR:HG21	4:A:793:HOH:O	1.87	0.74
1:B:75:LYS:HE2	4:B:852:HOH:O	1.89	0.73
1:A:307:LYS:H	1:A:311:ASN:HD21	1.35	0.72
1:B:180:ARG:NH1	1:B:248:VAL:O	2.23	0.71
1:C:143:GLN:H	1:C:144:PRO:HD3	1.55	0.70
1:C:115:HIS:HD2	1:C:226:PRO:HG2	1.57	0.70
1:A:143:GLN:N	1:A:144:PRO:CD	2.57	0.68
1:C:143:GLN:N	1:C:144:PRO:CD	2.57	0.68
1:B:132:PRO:CD	1:B:156:PHE:HD2	1.98	0.68
1:A:198:PHE:CE2	1:A:215:THR:HG22	2.29	0.68
1:B:211:VAL:O	1:B:215:THR:HG23	1.95	0.67
1:A:12:GLY:HA3	1:A:52:ASP:OD1	1.95	0.67
1:A:115:HIS:HD2	1:A:226:PRO:HG2	1.60	0.67
1:A:143:GLN:H	1:A:144:PRO:HD3	1.58	0.66
1:B:143:GLN:H	1:B:144:PRO:HD3	1.59	0.66
1:A:66:ILE:O	1:A:138:GLY:HA3	1.96	0.66
1:B:208:ILE:HB	4:B:705:HOH:O	1.95	0.66
1:B:143:GLN:N	1:B:144:PRO:CD	2.59	0.65
1:A:188:TYR:CD2	1:C:318:TYR:HB2	2.30	0.65
1:B:115:HIS:HD2	1:B:226:PRO:HG2	1.61	0.65
1:C:89:LYS:HD3	4:C:859:HOH:O	1.97	0.65
1:C:198:PHE:HE2	1:C:215:THR:HG22	1.62	0.65
1:A:187:HIS:O	1:A:191:LYS:HG2	1.97	0.64
1:C:187:HIS:O	1:C:191:LYS:HG2	1.97	0.64
4:B:763:HOH:O	1:C:201:THR:HG21	1.97	0.64
1:C:244:THR:C	1:C:282:LEU:HD12	2.22	0.64
1:B:15:PHE:O	1:B:99:GLY:HA3	1.99	0.63
1:B:135:THR:HB	1:B:137:SER:O	1.98	0.62
1:A:15:PHE:O	1:A:99:GLY:HA3	1.99	0.62
1:B:211:VAL:O	1:B:215:THR:CG2	2.47	0.62
1:A:96:VAL:HG23	4:A:718:HOH:O	2.00	0.62
1:A:308:ARG:O	1:B:180:ARG:HA	1.99	0.61
1:B:175:VAL:HG11	1:B:215:THR:HB	1.83	0.61
1:A:198:PHE:HE2	1:A:215:THR:HG22	1.64	0.61
1:B:152:LEU:HD13	1:B:193:LEU:HD21	1.83	0.60
1:C:66:ILE:O	1:C:138:GLY:HA3	2.02	0.60
1:C:152:LEU:HD13	1:C:193:LEU:HD21	1.83	0.60
1:B:233:VAL:HG13	1:B:241:THR:HB	1.83	0.60
1:B:187:HIS:O	1:B:191:LYS:HG2	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:12:GLY:HA3	1:C:52:ASP:OD1	2.03	0.59
1:A:152:LEU:HD13	1:A:193:LEU:HD21	1.83	0.59
1:A:254:TYR:CE2	1:A:258:LEU:HD11	2.38	0.59
1:A:311:ASN:O	1:B:184:PRO:HA	2.02	0.59
1:B:66:ILE:O	1:B:138:GLY:HA3	2.02	0.58
1:C:125:ALA:HA	1:C:178:GLY:O	2.03	0.58
1:C:15:PHE:O	1:C:99:GLY:HA3	2.02	0.58
1:A:233:VAL:HG13	1:A:241:THR:HB	1.84	0.57
1:A:180:ARG:HG3	1:A:248:VAL:HG11	1.86	0.57
1:C:233:VAL:HG13	1:C:241:THR:HB	1.85	0.57
1:B:12:GLY:HA3	1:B:52:ASP:OD1	2.05	0.56
1:A:6:LYS:HE2	1:A:91:GLY:O	2.06	0.56
1:A:115:HIS:CD2	1:A:226:PRO:HG2	2.41	0.56
1:C:115:HIS:CD2	1:C:226:PRO:HG2	2.40	0.55
1:C:180:ARG:NH2	1:C:235:GLY:O	2.40	0.55
1:A:127:THR:OG1	1:A:186:GLU:OE2	2.18	0.55
1:B:244:THR:HG22	1:B:277:GLU:O	2.06	0.55
1:A:156:PHE:HD1	1:A:156:PHE:O	1.89	0.54
1:B:78:GLU:HB2	1:B:164:TRP:CE2	2.41	0.54
1:A:156:PHE:O	1:A:156:PHE:CD1	2.60	0.54
1:C:180:ARG:CZ	1:C:251:GLY:HA2	2.38	0.54
1:C:317:ASP:OD2	1:C:319:LEU:OXT	2.25	0.54
1:A:283:GLY:HA3	1:A:289:VAL:HG23	1.90	0.53
1:A:115:HIS:HD2	1:A:226:PRO:CG	2.20	0.53
1:B:115:HIS:CD2	1:B:226:PRO:HG2	2.41	0.53
1:B:274:ASP:HB3	1:B:276:MET:SD	2.48	0.53
1:B:180:ARG:NH1	1:B:251:GLY:N	2.56	0.53
1:B:148:LEU:O	1:B:169:ILE:HD13	2.09	0.53
1:A:273:LEU:HD22	1:A:274:ASP:N	2.23	0.53
1:B:20:PRO:HD3	1:B:139:ASN:OD1	2.09	0.53
1:B:172:LYS:HD2	4:B:833:HOH:O	2.08	0.53
1:B:179:LEU:O	1:B:199:SER:HB2	2.09	0.53
1:A:77:ASN:OD1	1:A:103:MET:HA	2.10	0.52
1:A:318:TYR:CE2	1:B:184:PRO:HB2	2.44	0.52
1:B:180:ARG:NH1	1:B:251:GLY:HA3	2.23	0.52
1:C:253:SER:OG	1:C:256:GLU:HG3	2.09	0.52
1:A:148:LEU:O	1:A:169:ILE:HD13	2.09	0.52
1:C:115:HIS:HD2	1:C:226:PRO:CG	2.22	0.52
1:A:254:TYR:CZ	1:A:258:LEU:HD11	2.44	0.52
1:A:188:TYR:HE2	1:C:319:LEU:HG	1.75	0.52
1:A:118:LEU:C	1:A:118:LEU:HD12	2.34	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:180:ARG:HG3	1:C:248:VAL:HG11	1.92	0.51
1:A:132:PRO:CD	1:A:156:PHE:CD2	2.86	0.51
1:C:6:LYS:HE2	1:C:91:GLY:O	2.09	0.51
1:A:254:TYR:CE2	1:A:258:LEU:CD1	2.94	0.51
1:B:77:ASN:OD1	1:B:103:MET:HA	2.10	0.51
1:C:20:PRO:HD3	1:C:139:ASN:OD1	2.11	0.51
1:A:135:THR:HB	1:A:137:SER:O	2.10	0.51
1:B:115:HIS:HD2	1:B:226:PRO:CG	2.22	0.51
1:B:6:LYS:HE2	1:B:91:GLY:O	2.10	0.51
1:A:132:PRO:CD	1:A:156:PHE:HD2	2.08	0.51
1:C:77:ASN:OD1	1:C:103:MET:HA	2.11	0.50
1:A:312:HIS:HA	1:B:184:PRO:HB3	1.93	0.50
1:C:73:VAL:HG11	1:C:140:LEU:HD13	1.93	0.50
1:B:233:VAL:HG12	1:B:244:THR:HG21	1.93	0.50
1:C:156:PHE:O	1:C:156:PHE:CD1	2.64	0.50
1:C:20:PRO:HD3	1:C:139:ASN:CG	2.37	0.50
1:A:244:THR:HG22	1:A:277:GLU:O	2.12	0.50
1:A:89:LYS:NZ	4:A:773:HOH:O	2.45	0.49
1:B:118:LEU:HD12	1:B:118:LEU:C	2.36	0.49
1:C:118:LEU:C	1:C:118:LEU:HD12	2.37	0.49
1:B:245:GLY:HA2	1:B:282:LEU:CD1	2.43	0.49
1:A:253:SER:OG	1:A:256:GLU:HG3	2.12	0.49
1:C:273:LEU:HD22	1:C:274:ASP:N	2.28	0.49
1:C:148:LEU:O	1:C:169:ILE:HD13	2.13	0.49
1:A:246:THR:CG2	4:A:793:HOH:O	2.55	0.49
1:B:130:ASN:HB3	1:B:135:THR:HG23	1.94	0.48
1:C:135:THR:OG1	1:C:137:SER:O	2.31	0.48
1:B:7:PRO:HG2	4:B:739:HOH:O	2.13	0.48
1:B:180:ARG:NE	4:B:853:HOH:O	2.46	0.48
1:A:20:PRO:HD3	1:A:139:ASN:OD1	2.13	0.48
1:B:156:PHE:O	1:B:156:PHE:CD1	2.66	0.48
1:B:169:ILE:HD13	1:B:169:ILE:H	1.78	0.48
1:C:15:PHE:CZ	1:C:17:LYS:HB2	2.48	0.48
1:A:283:GLY:HA3	1:A:289:VAL:CG2	2.43	0.48
1:C:141:HIS:CD2	1:C:142:GLY:N	2.82	0.48
1:A:15:PHE:CZ	1:A:17:LYS:HB2	2.49	0.48
1:C:220:LEU:O	1:C:221:GLY:C	2.57	0.48
1:C:245:GLY:HA2	1:C:282:LEU:HD11	1.94	0.48
1:C:245:GLY:HA2	1:C:282:LEU:CD1	2.43	0.47
1:C:233:VAL:HG12	1:C:244:THR:HG21	1.96	0.47
1:B:78:GLU:HB2	1:B:164:TRP:NE1	2.29	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:GLY:HA2	1:A:282:LEU:CD1	2.45	0.47
1:B:73:VAL:HG11	1:B:140:LEU:HD13	1.96	0.47
1:A:180:ARG:NH2	1:A:235:GLY:O	2.47	0.47
4:A:819:HOH:O	1:B:201:THR:HG21	2.15	0.47
1:A:73:VAL:HG11	1:A:140:LEU:HD13	1.97	0.47
1:B:233:VAL:HG12	1:B:244:THR:CG2	2.45	0.47
1:A:119:CYS:HB2	1:A:219:LEU:HD22	1.98	0.46
1:C:244:THR:HG22	1:C:277:GLU:O	2.14	0.46
1:B:283:GLY:HA3	1:B:289:VAL:HG23	1.97	0.46
1:C:137:SER:C	1:C:139:ASN:H	2.23	0.46
1:A:6:LYS:HD2	1:A:7:PRO:HD3	1.97	0.46
1:B:198:PHE:CE2	1:B:215:THR:HG22	2.49	0.46
1:C:244:THR:O	1:C:282:LEU:HD12	2.15	0.46
1:C:21:ARG:HH11	1:C:282:LEU:CD1	2.29	0.46
1:B:132:PRO:CD	1:B:156:PHE:CD2	2.84	0.46
1:A:148:LEU:HB3	1:A:169:ILE:HD12	1.97	0.46
1:C:6:LYS:HD2	1:C:7:PRO:HD3	1.97	0.46
1:C:24:VAL:HG13	4:C:741:HOH:O	2.15	0.46
1:A:180:ARG:NH1	1:A:248:VAL:O	2.50	0.45
1:A:246:THR:OG1	3:A:1000:DIR:NH2	2.50	0.45
1:B:261:THR:HG22	1:B:303:CYS:SG	2.57	0.45
1:C:21:ARG:HH11	1:C:282:LEU:HD13	1.81	0.45
1:C:64:PHE:O	1:C:65:GLN:HB2	2.16	0.45
1:C:151:GLU:OE1	1:C:151:GLU:N	2.49	0.45
1:C:287:GLU:HG3	1:C:288:GLU:N	2.32	0.45
1:C:301:LEU:HD13	4:C:858:HOH:O	2.17	0.45
1:A:169:ILE:HD13	1:A:169:ILE:H	1.82	0.44
1:B:77:ASN:O	1:B:78:GLU:C	2.59	0.44
1:A:148:LEU:HB3	1:A:169:ILE:CD1	2.48	0.44
1:B:156:PHE:CD1	1:B:156:PHE:N	2.85	0.44
1:C:180:ARG:HD2	1:C:248:VAL:HG12	1.98	0.44
1:A:180:ARG:HD2	1:A:248:VAL:HG12	2.00	0.44
1:C:127:THR:OG1	1:C:186:GLU:OE2	2.26	0.44
1:A:64:PHE:O	1:A:65:GLN:HB2	2.17	0.44
1:C:156:PHE:CD1	1:C:156:PHE:N	2.84	0.44
1:A:38:GLU:CD	1:A:38:GLU:H	2.25	0.44
1:C:148:LEU:HB3	1:C:169:ILE:HD12	2.00	0.44
1:B:148:LEU:HB3	1:B:169:ILE:CD1	2.48	0.44
1:B:15:PHE:CZ	1:B:17:LYS:HB2	2.53	0.43
1:A:130:ASN:HB3	1:A:135:THR:HG23	2.00	0.43
1:A:180:ARG:CZ	1:A:251:GLY:HA2	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:148:LEU:HB3	1:C:169:ILE:CD1	2.48	0.43
1:C:246:THR:OG1	3:C:1002:DIR:NH2	2.51	0.43
1:B:180:ARG:NH1	1:B:251:GLY:HA2	2.29	0.43
1:B:283:GLY:HA3	1:B:289:VAL:CG2	2.48	0.43
1:C:198:PHE:CD2	1:C:215:THR:HG22	2.51	0.43
1:B:246:THR:HG21	3:B:1001:DIR:HH21	1.83	0.43
1:B:148:LEU:HB3	1:B:169:ILE:HD12	2.00	0.43
1:A:156:PHE:CD1	1:A:156:PHE:C	2.94	0.43
1:B:20:PRO:HD3	1:B:139:ASN:CG	2.44	0.43
1:B:289:VAL:O	1:B:293:VAL:HG23	2.19	0.43
1:A:233:VAL:HG12	1:A:244:THR:HG21	2.01	0.43
1:C:119:CYS:HB2	1:C:219:LEU:HD22	2.01	0.43
1:B:8:ILE:HD13	1:B:93:ILE:HB	2.01	0.43
1:B:64:PHE:O	1:B:65:GLN:HB2	2.19	0.42
1:A:38:GLU:CD	1:A:38:GLU:N	2.77	0.42
1:B:127:THR:HG23	1:B:179:LEU:HD22	2.01	0.42
1:B:180:ARG:CZ	4:B:853:HOH:O	2.67	0.42
1:B:180:ARG:HH12	1:B:251:GLY:HA3	1.84	0.42
1:C:169:ILE:HD13	1:C:169:ILE:H	1.82	0.42
1:B:126:HIS:CE1	3:B:1001:DIR:NH2	2.88	0.42
1:B:6:LYS:HD2	1:B:7:PRO:HD3	2.01	0.42
1:B:102:SER:HA	1:B:144:PRO:HG3	2.01	0.42
1:C:180:ARG:NH2	1:C:251:GLY:HA2	2.35	0.42
1:C:283:GLY:HA3	1:C:289:VAL:HG23	2.00	0.42
1:B:180:ARG:HH12	1:B:251:GLY:CA	2.30	0.41
1:C:233:VAL:HG13	1:C:241:THR:CB	2.50	0.41
1:A:313:LYS:HA	1:A:314:PRO:HD3	1.78	0.41
1:A:318:TYR:HE2	1:B:184:PRO:HB2	1.85	0.41
1:C:118:LEU:HD12	1:C:118:LEU:O	2.20	0.41
1:A:180:ARG:HD2	1:A:248:VAL:CG1	2.51	0.41
1:B:293:VAL:O	1:B:297:VAL:HG23	2.19	0.41
1:C:128:ASP:OD1	3:C:1002:DIR:NH1	2.53	0.41
1:C:214:GLU:O	1:C:215:THR:C	2.63	0.41
1:A:19:GLN:OE1	1:A:21:ARG:HB2	2.21	0.41
1:A:201:THR:HG22	1:C:308:ARG:HB2	2.02	0.41
1:B:246:THR:HG21	4:B:875:HOH:O	2.19	0.41
1:B:287:GLU:HG3	1:B:288:GLU:N	2.34	0.41
1:C:38:GLU:N	1:C:38:GLU:CD	2.79	0.41
1:A:287:GLU:HG3	1:A:288:GLU:N	2.36	0.41
1:B:166:THR:O	1:B:167:PRO:C	2.64	0.41
1:B:200:MET:HE1	1:B:252:LEU:HD23	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:VAL:HG12	1:A:244:THR:CG2	2.51	0.41
1:C:225:ARG:NH2	4:C:711:HOH:O	2.53	0.41
1:A:156:PHE:CD1	1:A:156:PHE:N	2.88	0.41
1:B:273:LEU:HD22	1:B:274:ASP:N	2.35	0.40
1:C:101:HIS:CE1	1:C:122:TRP:CZ2	3.09	0.40
1:B:10:ILE:N	1:B:10:ILE:HD12	2.36	0.40
1:B:233:VAL:HG13	1:B:241:THR:CB	2.51	0.40
1:C:86:GLU:O	1:C:89:LYS:HB3	2.22	0.40
1:C:180:ARG:CZ	1:C:251:GLY:CA	2.99	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	312/314 (99%)	300 (96%)	10 (3%)	2 (1%)	21	23
1	B	312/314 (99%)	300 (96%)	10 (3%)	2 (1%)	21	23
1	C	312/314 (99%)	297 (95%)	11 (4%)	4 (1%)	9	8
All	All	936/942 (99%)	897 (96%)	31 (3%)	8 (1%)	14	14

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	180	ARG
1	B	180	ARG
1	C	180	ARG
1	A	143	GLN
1	B	143	GLN
1	C	143	GLN
1	C	221	GLY
1	C	167	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	264/264 (100%)	246 (93%)	18 (7%)	14	17
1	B	264/264 (100%)	245 (93%)	19 (7%)	13	15
1	C	264/264 (100%)	246 (93%)	18 (7%)	14	17
All	All	792/792 (100%)	737 (93%)	55 (7%)	14	16

All (55) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	VAL
1	A	57	ASP
1	A	89	LYS
1	A	132	PRO
1	A	135	THR
1	A	152	LEU
1	A	156	PHE
1	A	167	PRO
1	A	169	ILE
1	A	201	THR
1	A	225	ARG
1	A	233	VAL
1	A	244	THR
1	A	246	THR
1	A	273	LEU
1	A	287	GLU
1	A	295	THR
1	A	313	LYS
1	B	24	VAL
1	B	57	ASP
1	B	89	LYS
1	B	132	PRO
1	B	135	THR
1	B	152	LEU
1	B	156	PHE
1	B	167	PRO

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Mol	Chain	Res	Type
1	B	169	ILE
1	B	201	THR
1	B	215	THR
1	B	225	ARG
1	B	233	VAL
1	B	244	THR
1	B	246	THR
1	B	273	LEU
1	B	287	GLU
1	B	295	THR
1	B	313	LYS
1	C	24	VAL
1	C	57	ASP
1	C	89	LYS
1	C	132	PRO
1	C	152	LEU
1	C	156	PHE
1	C	167	PRO
1	C	169	ILE
1	C	201	THR
1	C	222	ARG
1	C	233	VAL
1	C	236	LEU
1	C	244	THR
1	C	246	THR
1	C	273	LEU
1	C	287	GLU
1	C	295	THR
1	C	313	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	115	HIS
1	A	311	ASN
1	B	115	HIS
1	B	311	ASN
1	C	115	HIS
1	C	311	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](#)

Of 9 ligands modelled in this entry, 6 are monoatomic - leaving 3 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# $ Z > 2$	Counts	RMSZ	# $ Z > 2$
3	DIR	C	1002	2	8,10,10	2.72	4 (50%)	4,12,12	4.35	2 (50%)
3	DIR	B	1001	2	8,10,10	2.73	4 (50%)	4,12,12	4.53	2 (50%)
3	DIR	A	1000	2	8,10,10	2.61	4 (50%)	4,12,12	4.50	2 (50%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	DIR	C	1002	2	-	2/9/11/11	-
3	DIR	B	1001	2	-	2/9/11/11	-
3	DIR	A	1000	2	-	2/9/11/11	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	1000	DIR	O-C	4.84	1.36	1.22
3	C	1002	DIR	O-C	4.71	1.36	1.22
3	B	1001	DIR	O-C	4.71	1.36	1.22
3	C	1002	DIR	CD-NG	4.41	1.41	1.33
3	B	1001	DIR	CD-NG	4.20	1.41	1.33
3	A	1000	DIR	CD-NG	3.46	1.39	1.33
3	B	1001	DIR	OXT-C	-3.21	1.20	1.30
3	A	1000	DIR	OXT-C	-3.07	1.20	1.30
3	C	1002	DIR	OXT-C	-2.77	1.21	1.30
3	C	1002	DIR	OH1-NH1	-2.52	1.33	1.40
3	A	1000	DIR	OH1-NH1	-2.51	1.33	1.40
3	B	1001	DIR	OH1-NH1	-2.30	1.33	1.40

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1001	DIR	OH1-NH1-CD	8.13	119.77	109.97
3	A	1000	DIR	OH1-NH1-CD	7.94	119.54	109.97
3	C	1002	DIR	OH1-NH1-CD	7.63	119.17	109.97
3	C	1002	DIR	NH2-CD-NG	-3.91	112.90	119.17
3	A	1000	DIR	NH2-CD-NG	-3.85	112.99	119.17
3	B	1001	DIR	NH2-CD-NG	-3.62	113.35	119.17

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	1000	DIR	C-CA-CB-NG
3	B	1001	DIR	C-CA-CB-NG
3	C	1002	DIR	C-CA-CB-NG
3	A	1000	DIR	CA-CB-NG-CD
3	B	1001	DIR	CA-CB-NG-CD
3	C	1002	DIR	CA-CB-NG-CD

There are no ring outliers.

3 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	1002	DIR	2	0
3	B	1001	DIR	2	0
3	A	1000	DIR	1	0

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.