Full wwPDB X-ray Structure Validation Report (i)

Sep 20, 2023 – 09:36 PM EDT

PDB ID	:	5JO7
Title	:	Henbane premnaspirodiene synthase (HPS), also known as Henbane vetispiradiene synthase (HVS) from Hyoscyamus muticus
Authors	:	Koo, H.J.; Xu, Y.; Louie, G.V.; Bowman, M.; Noel, J.P.
Deposited on	:	2016-05-02
Resolution	:	2.15 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 (i) 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 (1)) were used in the production of this report:

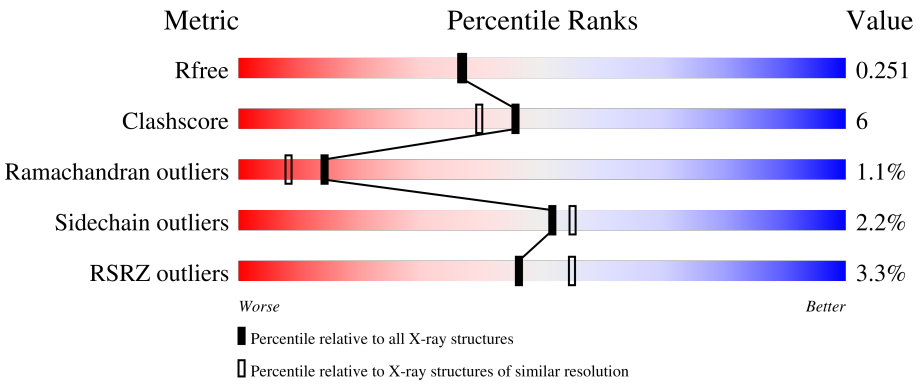
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.35.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 2.15 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}	130704	1479 (2.16-2.16)
Clashscore	141614	1585 (2.16-2.16)
Ramachandran outliers	138981	1560 (2.16-2.16)
Sidechain outliers	138945	1559 (2.16-2.16)
RSRZ outliers	127900	1456 (2.16-2.16)

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 <=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	538	2% 79% 15% • 5%
1	B	538	% 75% 17% • 6%
1	C	538	2% 79% 13% • 7%
1	D	538	7% 77% 15% • 7%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 17486 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 Vetispiradiene synthase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	510	4190	2683	690	800	17	0	1	0
1	B	506	4169	2674	686	792	17	0	2	0
1	C	503	4141	2658	677	789	17	0	2	0
1	D	499	4099	2629	672	781	17	0	1	0

There are 32 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	18	GLY	-	expression tag	UNP Q39978
A	19	SER	-	expression tag	UNP Q39978
A	20	HIS	-	expression tag	UNP Q39978
A	21	MET	-	expression tag	UNP Q39978
A	22	ALA	-	expression tag	UNP Q39978
A	30	ARG	HIS	conflict	UNP Q39978
A	381	ARG	GLY	conflict	UNP Q39978
A	387	ALA	GLY	conflict	UNP Q39978
B	18	GLY	-	expression tag	UNP Q39978
B	19	SER	-	expression tag	UNP Q39978
B	20	HIS	-	expression tag	UNP Q39978
B	21	MET	-	expression tag	UNP Q39978
B	22	ALA	-	expression tag	UNP Q39978
B	30	ARG	HIS	conflict	UNP Q39978
B	381	ARG	GLY	conflict	UNP Q39978
B	387	ALA	GLY	conflict	UNP Q39978
C	18	GLY	-	expression tag	UNP Q39978
C	19	SER	-	expression tag	UNP Q39978
C	20	HIS	-	expression tag	UNP Q39978
C	21	MET	-	expression tag	UNP Q39978
C	22	ALA	-	expression tag	UNP Q39978

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Chain	Residue	Modelled	Actual	Comment	Reference
C	30	ARG	HIS	conflict	UNP Q39978
C	381	ARG	GLY	conflict	UNP Q39978
C	387	ALA	GLY	conflict	UNP Q39978
D	18	GLY	-	expression tag	UNP Q39978
D	19	SER	-	expression tag	UNP Q39978
D	20	HIS	-	expression tag	UNP Q39978
D	21	MET	-	expression tag	UNP Q39978
D	22	ALA	-	expression tag	UNP Q39978
D	30	ARG	HIS	conflict	UNP Q39978
D	381	ARG	GLY	conflict	UNP Q39978
D	387	ALA	GLY	conflict	UNP Q39978

- Molecule 2 is water.

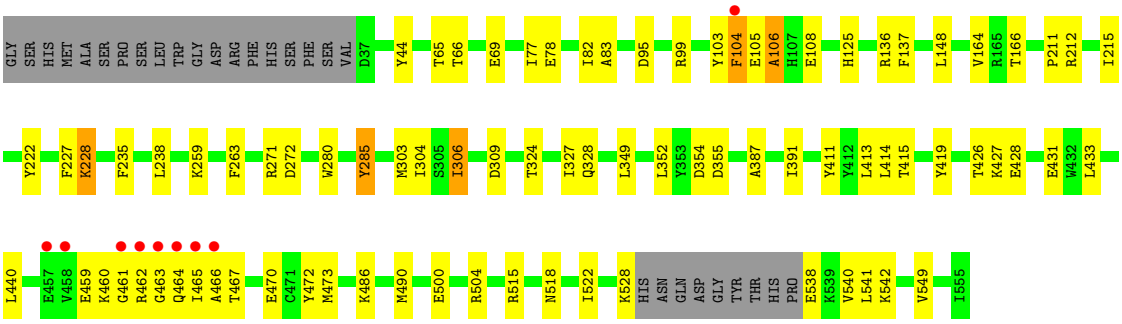
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	306	306	306	0	0
			Total	O		
2	B	252	252	252	0	0
			Total	O		
2	C	198	198	198	0	0
			Total	O		
2	D	131	131	131	0	0

3 Residue-property plots (i)

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.

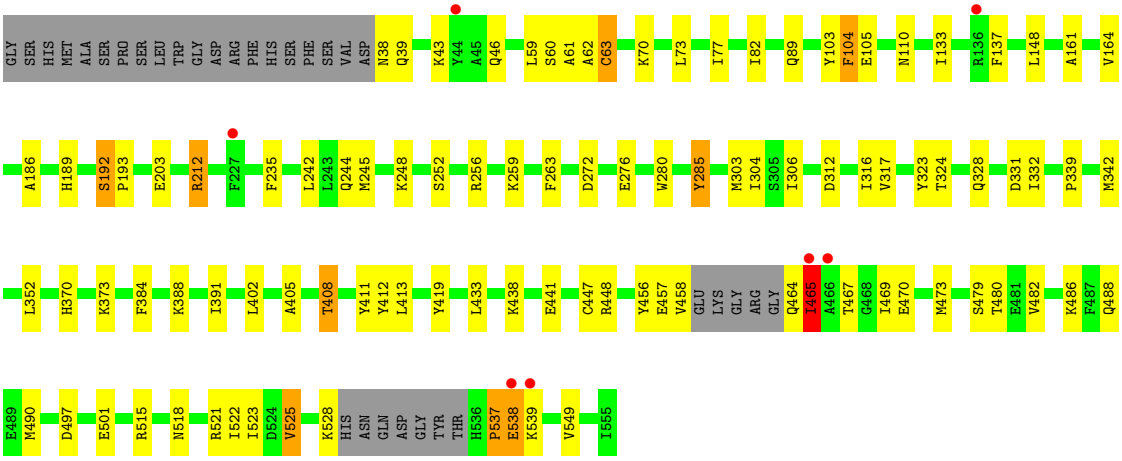
• Molecule 1: Vetispiradiene synthase 1

Chain A:



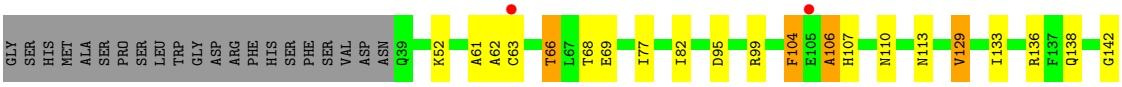
• Molecule 1: Vetispiradiene synthase 1

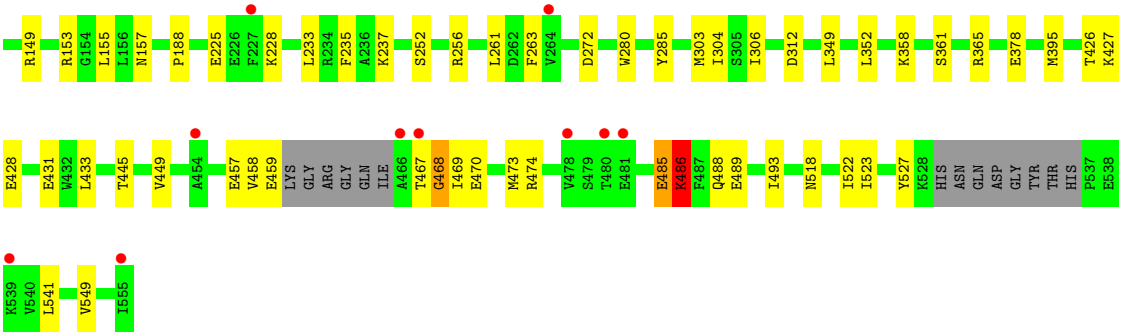
Chain B:



• Molecule 1: Vetispiradiene synthase 1

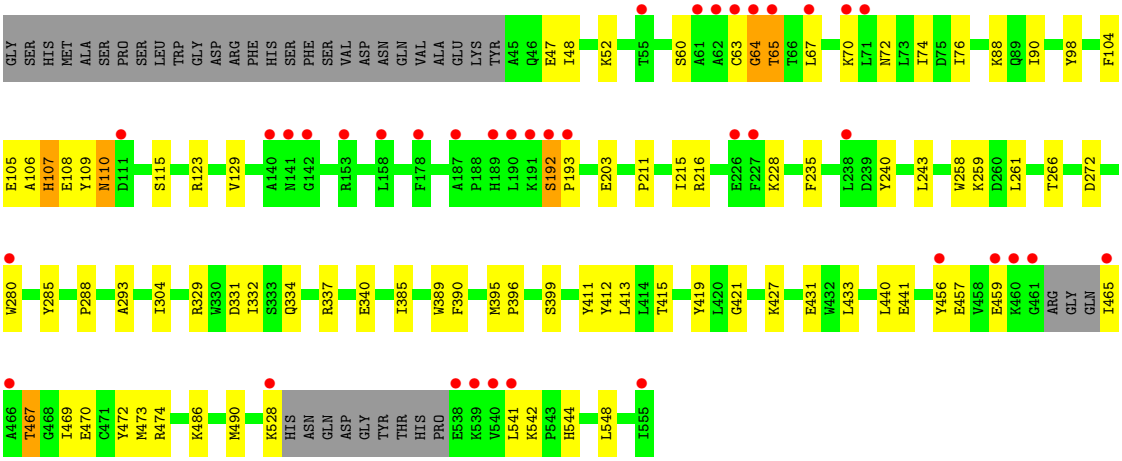
Chain C:





- Molecule 1: Vetispiradiene synthase 1

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	100.45Å 100.69Å 222.73Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	74.69 – 2.15 74.69 – 2.15	Depositor EDS
% Data completeness (in resolution range)	85.0 (74.69-2.15) 84.7 (74.69-2.15)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.88 (at 2.14Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.191 , 0.250 0.193 , 0.251	Depositor DCC
R_{free} test set	2192 reflections (2.09%)	wwPDB-VP
Wilson B-factor (Å^2)	24.6	Xtriage
Anisotropy	0.591	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 49.4	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.016 for k,h,-l	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	17486	wwPDB-VP
Average B, all atoms (Å^2)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 14.08% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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	0.47	0/4283	0.60	1/5800 (0.0%)
1	B	0.40	0/4267	0.57	0/5781
1	C	0.40	0/4237	0.60	4/5739 (0.1%)
1	D	0.36	0/4190	0.52	0/5674
All	All	0.41	0/16977	0.57	5/22994 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	B	0	2
1	C	0	1
1	D	0	1
All	All	0	5

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	485	GLU	C-N-CA	6.60	138.20	121.70
1	A	349	LEU	CA-CB-CG	6.24	129.65	115.30
1	C	485	GLU	N-CA-C	5.98	127.14	111.00
1	C	468	GLY	N-CA-C	-5.27	99.92	113.10
1	C	485	GLU	CA-C-N	5.24	128.72	117.20

There are no chirality outliers.

All (5) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	106	ALA	Peptide
1	B	104	PHE	Peptide
1	B	59	LEU	Peptide
1	C	106	ALA	Peptide
1	D	64	GLY	Peptide

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	4190	0	4135	50	0
1	B	4169	0	4113	68	0
1	C	4141	0	4092	48	0
1	D	4099	0	4050	50	0
2	A	306	0	0	3	0
2	B	252	0	0	7	0
2	C	198	0	0	6	0
2	D	131	0	0	3	0
All	All	17486	0	16390	212	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 6.

All (212) 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:464:GLN:NE2	1:B:465:ILE:O	2.09	0.85
1:A:271:ARG:NH1	1:A:309:ASP:OD1	2.11	0.84
1:D:203:GLU:O	1:D:528:LYS:NZ	2.16	0.79
1:D:340:GLU:OE1	2:D:601:HOH:O	2.03	0.77
1:C:395:MET:HG2	1:C:474:ARG:HD3	1.68	0.76
1:B:61:ALA:HB1	1:B:63:CYS:H	1.51	0.75
1:D:63:CYS:SG	1:D:70:LYS:NZ	2.59	0.74
1:C:225:GLU:O	1:C:228:LYS:NZ	2.19	0.72
1:B:61:ALA:HA	1:B:62:ALA:HB3	1.72	0.70
1:A:465:ILE:O	1:A:467:THR:N	2.22	0.69
1:D:395:MET:HG2	1:D:474:ARG:HD3	1.74	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:104:PHE:H	1:A:104:PHE:HD1	1.42	0.67
1:A:271:ARG:HG3	1:A:271:ARG:HH11	1.59	0.67
1:B:486:LYS:NZ	2:B:601:HOH:O	2.19	0.67
1:A:104:PHE:O	1:A:136:ARG:NH2	2.24	0.66
1:B:408:THR:HG21	2:B:628:HOH:O	1.96	0.65
1:D:64:GLY:H	1:D:65:THR:HA	1.61	0.65
1:B:537:PRO:O	1:B:539:LYS:N	2.30	0.64
1:A:103:TYR:O	1:A:105:GLU:N	2.31	0.63
1:B:339:PRO:HD2	1:B:342:MET:HE2	1.81	0.63
1:A:472:TYR:CZ	1:A:486:LYS:HG2	2.32	0.63
1:C:153:ARG:NH2	2:C:606:HOH:O	2.30	0.62
1:D:192:SER:OG	1:D:193:PRO:HD2	2.01	0.61
1:C:358:LYS:NZ	2:C:608:HOH:O	2.32	0.60
1:B:192:SER:OG	1:B:193:PRO:HD2	2.00	0.60
1:B:285:TYR:CD2	1:B:515:ARG:HD2	2.36	0.60
1:B:61:ALA:HB1	1:B:63:CYS:N	2.17	0.60
1:D:243:LEU:HD21	1:D:288:PRO:HG3	1.82	0.59
1:B:456:TYR:C	1:B:458:VAL:H	2.06	0.59
1:B:60:SER:N	1:B:61:ALA:HB2	2.17	0.58
1:C:188:PRO:O	2:C:601:HOH:O	2.17	0.58
1:C:66:THR:HG23	1:C:68:THR:H	1.68	0.58
1:B:408:THR:HG22	1:B:448:ARG:NH2	2.18	0.58
1:D:74:ILE:HD12	1:D:90:ILE:HG23	1.86	0.58
1:D:456:TYR:HD1	1:D:467:THR:HG21	1.68	0.57
1:A:303:MET:HG3	1:A:352:LEU:HD22	1.87	0.57
1:D:109:TYR:O	1:D:115:SER:OG	2.20	0.57
1:D:334:GLN:OE1	1:D:337:ARG:NH1	2.34	0.56
1:B:304:ILE:HD11	1:B:413:LEU:HD23	1.88	0.56
1:A:470:GLU:HA	1:A:473:MET:HE3	1.86	0.56
1:A:538:GLU:HB2	1:A:540:VAL:HG23	1.88	0.56
1:D:98:TYR:CE1	1:D:129:VAL:HG12	2.41	0.56
1:B:408:THR:HG22	1:B:448:ARG:HH21	1.70	0.56
1:B:323:TYR:CD1	1:B:342:MET:HE3	2.41	0.56
1:B:303:MET:HG3	1:B:352:LEU:HD22	1.87	0.55
1:D:64:GLY:N	1:D:65:THR:HA	2.21	0.55
1:A:137:PHE:HB3	1:A:148:LEU:HD11	1.88	0.55
1:A:285:TYR:CD2	1:A:515:ARG:HD2	2.41	0.55
1:C:52:LYS:NZ	2:C:614:HOH:O	2.40	0.54
1:B:203:GLU:O	1:B:528:LYS:NZ	2.40	0.54
1:C:106:ALA:CB	1:C:107:HIS:HA	2.37	0.54
1:C:518:ASN:O	1:C:522[B]:ILE:HG13	2.07	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:125:HIS:O	1:A:504:ARG:NH1	2.37	0.54
1:A:78:GLU:HA	1:A:83:ALA:HB2	1.89	0.54
1:A:515:ARG:NH2	2:A:619:HOH:O	2.40	0.54
1:C:104:PHE:CG	1:C:133:ILE:HG13	2.43	0.54
1:C:470:GLU:HA	1:C:473:MET:HE3	1.90	0.54
1:B:467:THR:HG23	1:B:470:GLU:H	1.73	0.53
1:A:66:THR:N	1:A:69:GLU:OE2	2.41	0.53
1:B:419:TYR:CE2	1:B:433:LEU:HD11	2.44	0.53
1:D:258:TRP:NE1	1:D:272:ASP:OD1	2.40	0.53
1:B:518:ASN:O	1:B:522:ILE:HG13	2.09	0.53
1:A:263:PHE:CE2	1:A:306:ILE:HG13	2.45	0.52
1:D:419:TYR:CE2	1:D:433:LEU:HD11	2.44	0.52
1:D:472:TYR:CZ	1:D:486:LYS:HG2	2.45	0.52
1:A:463:GLY:HA3	1:A:464:GLN:C	2.29	0.51
1:A:95:ASP:O	1:A:99:ARG:HG2	2.11	0.51
1:B:469:ILE:O	1:B:473:MET:HG3	2.11	0.51
1:C:77:ILE:HG23	1:C:82:ILE:HG13	1.91	0.51
1:C:426:THR:OG1	1:C:428:GLU:HG2	2.11	0.51
1:B:456:TYR:O	1:B:458:VAL:N	2.43	0.51
1:C:303:MET:HG3	1:C:352:LEU:HD22	1.92	0.51
1:A:486:LYS:O	1:A:490:MET:HG3	2.11	0.51
1:C:106:ALA:HB1	1:C:107:HIS:HA	1.92	0.51
1:A:472:TYR:CE2	1:A:486:LYS:HG2	2.46	0.50
1:C:485:GLU:N	1:C:486:LYS:HB3	2.26	0.50
1:A:65:THR:HB	1:A:69:GLU:HG3	1.94	0.50
1:B:137:PHE:HB3	1:B:148:LEU:HD11	1.94	0.50
1:C:110:ASN:OD1	1:C:136:ARG:HD3	2.12	0.50
1:C:61:ALA:C	1:C:63:CYS:H	2.15	0.50
1:C:104:PHE:HZ	1:C:129:VAL:HG23	1.77	0.50
1:C:113:ASN:OD1	1:C:157:ASN:ND2	2.38	0.49
1:A:212:ARG:HD2	2:A:744:HOH:O	2.12	0.49
1:D:123:ARG:HG2	1:D:129:VAL:HG22	1.93	0.49
1:A:166:THR:OG1	1:A:500:GLU:HB2	2.13	0.49
1:C:427:LYS:O	1:C:431:GLU:HG2	2.12	0.49
1:A:280[B]:TRP:HZ3	1:A:541:LEU:HD11	1.78	0.49
1:B:497:ASP:O	1:B:501:GLU:HG2	2.13	0.49
1:C:280[B]:TRP:HZ3	1:C:541:LEU:HD11	1.78	0.49
1:D:469:ILE:O	1:D:473:MET:HG3	2.12	0.49
1:B:285:TYR:CE2	1:B:515:ARG:HD2	2.48	0.49
1:C:469:ILE:O	1:C:473:MET:HG3	2.13	0.49
1:D:48:ILE:HG23	1:D:235:PHE:CE1	2.48	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:60:SER:OG	1:B:61:ALA:HA	2.14	0.48
1:D:107:HIS:HA	1:D:109:TYR:N	2.28	0.48
1:B:212:ARG:HD2	2:B:770:HOH:O	2.14	0.48
1:A:463:GLY:HA3	1:A:464:GLN:O	2.14	0.48
1:C:485:GLU:O	1:C:488:GLN:HB2	2.14	0.48
1:D:304:ILE:HD11	1:D:413:LEU:HD23	1.95	0.48
1:A:272:ASP:HB3	1:B:272:ASP:HB3	1.96	0.48
1:C:489:GLU:O	1:C:493:ILE:HG12	2.14	0.47
1:C:233:LEU:HG	1:C:237:LYS:HE2	1.96	0.47
1:D:329:ARG:NE	2:D:604:HOH:O	2.29	0.47
1:B:323:TYR:HB2	1:B:342:MET:CE	2.45	0.47
1:B:323:TYR:HD1	1:B:342:MET:HE3	1.79	0.47
1:C:138:GLN:NE2	1:C:142:GLY:O	2.48	0.47
1:C:272:ASP:HB3	1:D:272:ASP:HB3	1.97	0.47
1:C:378:GLU:OE2	2:C:602:HOH:O	2.20	0.47
1:B:263:PHE:CE2	1:B:306:ILE:HG13	2.49	0.47
1:A:66:THR:O	1:A:69:GLU:HG2	2.15	0.47
1:B:479:SER:OG	1:B:482:VAL:HG12	2.15	0.47
1:A:324:THR:O	1:A:328:GLN:HG2	2.15	0.47
1:B:252:SER:OG	1:B:256:ARG:NH1	2.48	0.47
1:D:211:PRO:O	1:D:215:ILE:HG12	2.15	0.47
1:A:355:ASP:OD1	2:A:601:HOH:O	2.20	0.46
1:B:259:LYS:HD3	1:B:259:LYS:HA	1.69	0.46
1:D:70:LYS:HA	1:D:70:LYS:HD3	1.58	0.46
1:D:107:HIS:HB2	1:D:110:ASN:OD1	2.15	0.46
1:A:542:LYS:NZ	1:B:312:ASP:OD2	2.48	0.46
1:B:60:SER:HA	1:B:89:GLN:HE21	1.81	0.46
1:C:467:THR:OG1	1:C:468:GLY:HA3	2.15	0.46
1:D:389:TRP:CZ3	1:D:396:PRO:HG3	2.51	0.46
1:B:384:PHE:CZ	1:B:388:LYS:HD2	2.51	0.46
1:D:427:LYS:O	1:D:431:GLU:HG2	2.16	0.46
1:B:61:ALA:HB3	1:B:70:LYS:NZ	2.31	0.45
1:C:303:MET:HE2	1:C:349:LEU:HD13	1.99	0.45
1:B:317:VAL:CG2	1:B:391:ILE:HG13	2.47	0.45
1:B:521:ARG:O	1:B:525:VAL:HG13	2.16	0.45
1:C:252:SER:O	1:C:256:ARG:HG3	2.16	0.45
1:A:304:ILE:HD11	1:A:413:LEU:HD23	1.98	0.45
1:C:486:LYS:N	1:C:489:GLU:H	2.15	0.45
1:B:324:THR:O	1:B:328:GLN:HG2	2.16	0.45
1:B:488:GLN:NE2	2:B:624:HOH:O	2.49	0.45
1:D:331:ASP:OD1	1:D:332:ILE:N	2.50	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:43:LYS:NZ	2:B:609:HOH:O	2.37	0.44
1:B:104:PHE:HD2	1:B:133:ILE:O	2.00	0.44
1:B:70:LYS:HD3	1:B:70:LYS:HA	1.74	0.44
1:C:95:ASP:O	1:C:99:ARG:HG2	2.17	0.44
1:C:104:PHE:O	1:C:136:ARG:NH2	2.47	0.44
1:C:149:ARG:HA	1:C:155:LEU:HD11	1.99	0.44
1:D:107:HIS:HA	1:D:108:GLU:C	2.37	0.44
1:A:427:LYS:O	1:A:431:GLU:HG2	2.17	0.44
1:B:161:ALA:O	1:B:164:VAL:HG22	2.18	0.44
1:C:445:THR:O	1:C:449:VAL:HB	2.17	0.44
1:C:395:MET:HG2	1:C:474:ARG:CD	2.44	0.44
1:A:459:GLU:O	1:A:461:GLY:N	2.51	0.44
1:B:38:ASN:HA	1:B:39:GLN:HA	1.82	0.44
1:A:518:ASN:O	1:A:522:ILE:HG13	2.18	0.43
1:B:373:LYS:NZ	2:B:612:HOH:O	2.44	0.43
1:C:365:ARG:HD3	2:C:702:HOH:O	2.17	0.43
1:C:523:ILE:HD11	1:C:527:TYR:HE2	1.83	0.43
1:A:77:ILE:HG23	1:A:82:ILE:HG13	2.01	0.43
1:A:105:GLU:O	1:A:106:ALA:HB3	2.18	0.43
1:B:331:ASP:OD1	1:B:332:ILE:N	2.52	0.43
1:B:339:PRO:HD2	1:B:342:MET:CE	2.48	0.43
1:C:228:LYS:HD3	1:C:233:LEU:HD22	2.00	0.43
1:A:65:THR:HB	1:A:69:GLU:CG	2.49	0.43
1:C:486:LYS:H	1:C:489:GLU:CB	2.31	0.43
1:B:186:ALA:HA	1:B:189[A]:HIS:NE2	2.34	0.43
1:C:263:PHE:CE2	1:C:306:ILE:HG13	2.53	0.43
1:D:105:GLU:HA	1:D:106:ALA:HA	1.90	0.43
1:D:329:ARG:HD3	2:D:608:HOH:O	2.18	0.43
1:A:222:TYR:CZ	1:A:228:LYS:HB2	2.54	0.43
1:A:415:THR:HG21	1:A:440:LEU:HD13	2.01	0.42
1:B:316:ILE:N	2:B:621:HOH:O	2.48	0.42
1:B:280[A]:TRP:HZ3	1:B:411:TYR:OH	2.03	0.42
1:D:104:PHE:N	1:D:104:PHE:CD2	2.87	0.42
1:A:528:LYS:HD2	1:A:528:LYS:HA	1.83	0.42
1:B:412:TYR:OH	1:B:441:GLU:OE1	2.23	0.42
1:D:293:ALA:HA	1:D:421:GLY:HA3	2.01	0.42
1:D:390:PHE:CD1	1:D:465:ILE:HG23	2.54	0.42
1:D:259:LYS:HA	1:D:259:LYS:HD3	1.82	0.42
1:A:263:PHE:CD2	1:A:306:ILE:HG13	2.53	0.42
1:D:415:THR:HG21	1:D:440:LEU:HD13	2.00	0.42
1:A:419:TYR:CE2	1:A:433:LEU:HD11	2.54	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:104:PHE:N	1:D:104:PHE:HD2	2.18	0.42
1:D:243:LEU:HD23	1:D:243:LEU:HA	1.65	0.42
1:D:486:LYS:O	1:D:490:MET:HG3	2.20	0.42
1:B:447:CYS:HA	1:B:523:ILE:HG12	2.02	0.41
1:D:216:ARG:HD3	1:D:240:TYR:CZ	2.55	0.41
1:D:261:LEU:HD23	1:D:261:LEU:HA	1.89	0.41
1:D:72:ASN:O	1:D:76:ILE:HG12	2.19	0.41
1:D:544:HIS:O	1:D:548:LEU:HB2	2.20	0.41
1:B:60:SER:HA	1:B:89:GLN:NE2	2.35	0.41
1:B:248:LYS:HE2	1:B:549:VAL:O	2.21	0.41
1:C:66:THR:O	1:C:69:GLU:HG2	2.20	0.41
1:D:412:TYR:OH	1:D:441:GLU:OE1	2.33	0.41
1:B:103:TYR:HB3	1:B:105:GLU:HG3	2.02	0.41
1:B:73:LEU:O	1:B:77:ILE:HG12	2.19	0.41
1:D:470:GLU:HA	1:D:473:MET:HE3	2.03	0.41
1:B:242:LEU:HA	1:B:245:MET:HE2	2.02	0.41
1:B:276:GLU:OE2	1:B:538:GLU:HG3	2.21	0.41
1:A:44:TYR:CD1	1:A:238:LEU:HD22	2.56	0.41
1:A:306:ILE:HD13	1:A:306:ILE:HA	1.93	0.41
1:B:43:LYS:O	1:B:46:GLN:HB2	2.21	0.41
1:B:405:ALA:HA	1:B:408:THR:HB	2.02	0.41
1:D:280[B]:TRP:HZ3	1:D:541:LEU:HD11	1.85	0.41
1:D:385:ILE:HA	1:D:385:ILE:HD13	1.85	0.41
1:B:77:ILE:HG23	1:B:82:ILE:HG13	2.03	0.41
1:B:104:PHE:HZ	1:B:110:ASN:HA	1.85	0.41
1:B:486:LYS:O	1:B:490:MET:HG3	2.21	0.41
1:A:426:THR:OG1	1:A:428:GLU:HG2	2.21	0.40
1:A:470:GLU:HA	1:A:473:MET:CE	2.52	0.40
1:B:244:GLN:HG2	1:B:248:LYS:HE3	2.03	0.40
1:D:52:LYS:HG2	1:D:235:PHE:CZ	2.56	0.40
1:A:211:PRO:O	1:A:215:ILE:HG12	2.21	0.40
1:A:259:LYS:NZ	1:A:272:ASP:OD2	2.55	0.40
1:C:459:GLU:OE2	1:C:459:GLU:N	2.44	0.40
1:C:312:ASP:OD2	1:D:542:LYS:NZ	2.54	0.40
1:A:387:ALA:O	1:A:391:ILE:HG12	2.22	0.40
1:C:261:LEU:HD23	1:C:261:LEU:HA	1.93	0.40
1:C:304:ILE:HD12	1:C:304:ILE:HG23	1.88	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	507/538 (94%)	489 (96%)	14 (3%)	4 (1%)	19 12
1	B	502/538 (93%)	485 (97%)	12 (2%)	5 (1%)	15 9
1	C	499/538 (93%)	479 (96%)	15 (3%)	5 (1%)	15 9
1	D	494/538 (92%)	471 (95%)	15 (3%)	8 (2%)	9 4
All	All	2002/2152 (93%)	1924 (96%)	56 (3%)	22 (1%)	14 8

All (22) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	104	PHE
1	A	460	LYS
1	A	462	ARG
1	A	466	ALA
1	B	192	SER
1	B	465	ILE
1	C	62	ALA
1	C	104	PHE
1	C	457	GLU
1	C	458	VAL
1	C	486	LYS
1	D	47	GLU
1	D	60	SER
1	D	467	THR
1	B	457	GLU
1	B	537	PRO
1	B	538	GLU
1	D	110	ASN
1	D	192	SER
1	D	457	GLU
1	D	459	GLU
1	D	107	HIS

5.3.2 Protein sidechains (i)

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	458/481 (95%)	446 (97%)	12 (3%)	46	47
1	B	457/481 (95%)	446 (98%)	11 (2%)	49	51
1	C	454/481 (94%)	446 (98%)	8 (2%)	59	63
1	D	449/481 (93%)	441 (98%)	8 (2%)	59	63
All	All	1818/1924 (94%)	1779 (98%)	39 (2%)	52	57

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	108	GLU
1	A	164	VAL
1	A	227	PHE
1	A	228	LYS
1	A	235	PHE
1	A	285	TYR
1	A	306	ILE
1	A	327	ILE
1	A	354	ASP
1	A	411	TYR
1	A	414	LEU
1	A	549	VAL
1	B	63	CYS
1	B	212	ARG
1	B	235	PHE
1	B	285	TYR
1	B	370	HIS
1	B	402	LEU
1	B	408	THR
1	B	438	LYS
1	B	465	ILE
1	B	480	THR
1	B	525	VAL
1	C	66	THR

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Mol	Chain	Res	Type
1	C	129	VAL
1	C	235	PHE
1	C	285	TYR
1	C	361	SER
1	C	433	LEU
1	C	486	LYS
1	C	549	VAL
1	D	65	THR
1	D	67	LEU
1	D	88	LYS
1	D	228	LYS
1	D	266	THR
1	D	285	TYR
1	D	399	SER
1	D	411	TYR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 monosaccharides 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	510/538 (94%)	-0.23	9 (1%)	68	75	12, 26, 58, 133	0
1	B	506/538 (94%)	-0.22	7 (1%)	75	80	17, 33, 65, 116	0
1	C	503/538 (93%)	-0.11	12 (2%)	59	67	19, 38, 76, 106	0
1	D	499/538 (92%)	0.27	38 (7%)	13	19	21, 52, 95, 126	0
All	All	2018/2152 (93%)	-0.08	66 (3%)	46	55	12, 37, 81, 133	0

All (66) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	65	THR	7.4
1	D	227	PHE	6.8
1	C	227	PHE	6.5
1	C	63	CYS	5.0
1	B	44	TYR	4.6
1	D	555	ILE	4.6
1	A	104	PHE	4.4
1	B	465	ILE	4.3
1	A	466	ALA	4.2
1	A	465	ILE	4.2
1	B	538	GLU	4.0
1	A	462	ARG	3.9
1	D	64	GLY	3.9
1	B	227	PHE	3.9
1	D	63	CYS	3.8
1	D	158	LEU	3.7
1	D	141	ASN	3.7
1	D	61	ALA	3.4
1	D	465	ILE	3.4
1	D	461	GLY	3.3
1	C	467	THR	3.3

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Mol	Chain	Res	Type	RSRZ
1	B	539	LYS	3.2
1	D	142	GLY	3.2
1	D	541	LEU	3.1
1	D	153	ARG	3.0
1	D	178	PHE	3.0
1	D	459	GLU	2.9
1	A	457	GLU	2.8
1	D	111	ASP	2.7
1	A	458	VAL	2.7
1	C	478	VAL	2.7
1	D	71	LEU	2.7
1	D	140	ALA	2.6
1	D	190	LEU	2.6
1	D	189	HIS	2.5
1	C	539	LYS	2.5
1	D	280[A]	TRP	2.5
1	D	187	ALA	2.5
1	C	555	ILE	2.5
1	D	460	LYS	2.5
1	D	466	ALA	2.5
1	D	528	LYS	2.4
1	D	538	GLU	2.4
1	D	456	TYR	2.4
1	B	466	ALA	2.4
1	D	192	SER	2.4
1	D	67	LEU	2.3
1	D	70	LYS	2.3
1	A	461	GLY	2.3
1	D	539	LYS	2.3
1	C	481	GLU	2.3
1	A	464	GLN	2.3
1	C	466	ALA	2.3
1	C	264	VAL	2.3
1	D	193	PRO	2.3
1	C	480	THR	2.2
1	D	540	VAL	2.2
1	C	105	GLU	2.2
1	D	238	LEU	2.2
1	D	191	LYS	2.2
1	D	62	ALA	2.2
1	D	226	GLU	2.1
1	B	136	ARG	2.1

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Mol	Chain	Res	Type	RSRZ
1	D	55	THR	2.1
1	C	454	ALA	2.0
1	A	463	GLY	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides 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.