



Full wwPDB X-ray Structure Validation Report

Feb 28, 2014 – 10:55 PM GMT

PDB ID : 3TBG
Title : Human cytochrome P450 2D6 with two thioridazines bound in active site
Authors : Wang, A.; Stout, C.D.; Johnson, E.F.
Deposited on : 2011-08-05
Resolution : 2.10 Å(reported)

This is a full wwPDB validation report for a publicly released PDB entry.
We welcome your comments at validation@mail.wwpdb.org
A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

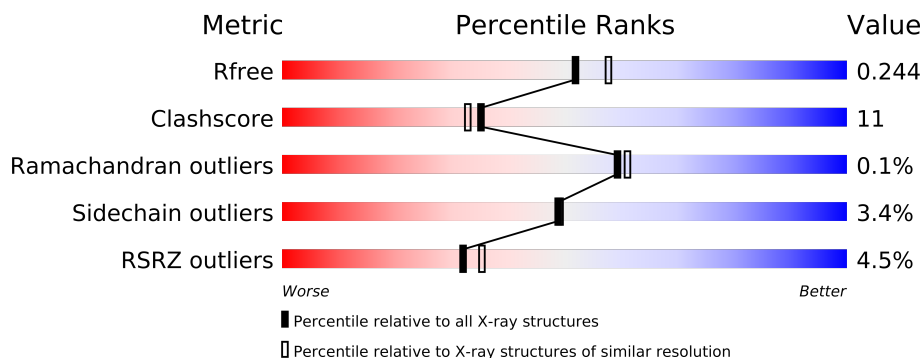
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.15 2013
Xtriage (Phenix) : dev-1323
EDS : stable22639
Percentile statistics : 21963
Refmac : 5.8.0049
CCP4 : 6.3.0 (Settle)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP) : stable22683

1 Overall quality at a glance

The reported resolution of this entry is 2.10 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}	66092	3012 (2.10-2.10)
Clashscore	79885	3649 (2.10-2.10)
Ramachandran outliers	78287	3610 (2.10-2.10)
Sidechain outliers	78261	3611 (2.10-2.10)
RSRZ outliers	66119	3013 (2.10-2.10)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	479	
1	B	479	
1	C	479	
1	D	479	

The following table lists non-polymeric compounds that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Geometry	Electron density
4	GOL	A	751	-	X
4	GOL	A	752	-	X
4	GOL	A	753	-	X
4	GOL	B	750	-	X
4	GOL	D	750	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 15538 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Cytochrome P450 2D6.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	467	Total	C	N	O	S	0	0	0
			3690	2367	654	655	14			
1	B	456	Total	C	N	O	S	0	0	0
			3614	2317	640	643	14			
1	C	467	Total	C	N	O	S	0	0	0
			3690	2367	654	655	14			
1	D	456	Total	C	N	O	S	0	0	0
			3614	2317	640	643	14			

There are 60 discrepancies between the modelled and reference sequences:

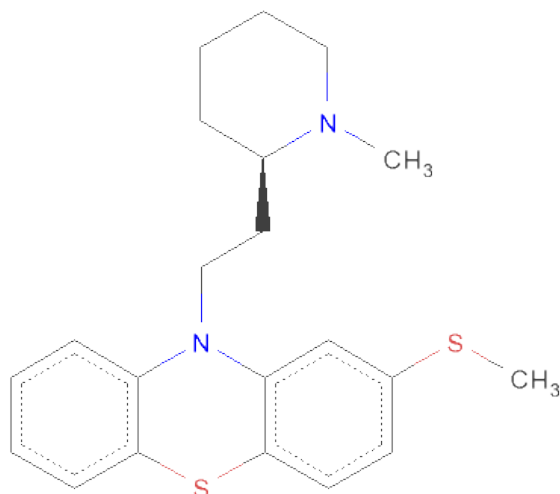
Chain	Residue	Modelled	Actual	Comment	Reference
A	23	MET	-	EXPRESSION TAG	UNP P10635
A	24	ALA	-	EXPRESSION TAG	UNP P10635
A	25	LYS	-	EXPRESSION TAG	UNP P10635
A	26	LYS	-	EXPRESSION TAG	UNP P10635
A	27	THR	-	EXPRESSION TAG	UNP P10635
A	28	SER	-	EXPRESSION TAG	UNP P10635
A	29	SER	-	EXPRESSION TAG	UNP P10635
A	30	LYS	-	EXPRESSION TAG	UNP P10635
A	31	GLY	-	EXPRESSION TAG	UNP P10635
A	32	LYS	-	EXPRESSION TAG	UNP P10635
A	33	LEU	-	EXPRESSION TAG	UNP P10635
A	498	HIS	-	EXPRESSION TAG	UNP P10635
A	499	HIS	-	EXPRESSION TAG	UNP P10635
A	500	HIS	-	EXPRESSION TAG	UNP P10635
A	501	HIS	-	EXPRESSION TAG	UNP P10635
B	23	MET	-	EXPRESSION TAG	UNP P10635
B	24	ALA	-	EXPRESSION TAG	UNP P10635
B	25	LYS	-	EXPRESSION TAG	UNP P10635
B	26	LYS	-	EXPRESSION TAG	UNP P10635
B	27	THR	-	EXPRESSION TAG	UNP P10635
B	28	SER	-	EXPRESSION TAG	UNP P10635

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Chain	Residue	Modelled	Actual	Comment	Reference
B	29	SER	-	EXPRESSION TAG	UNP P10635
B	30	LYS	-	EXPRESSION TAG	UNP P10635
B	31	GLY	-	EXPRESSION TAG	UNP P10635
B	32	LYS	-	EXPRESSION TAG	UNP P10635
B	33	LEU	-	EXPRESSION TAG	UNP P10635
B	498	HIS	-	EXPRESSION TAG	UNP P10635
B	499	HIS	-	EXPRESSION TAG	UNP P10635
B	500	HIS	-	EXPRESSION TAG	UNP P10635
B	501	HIS	-	EXPRESSION TAG	UNP P10635
C	23	MET	-	EXPRESSION TAG	UNP P10635
C	24	ALA	-	EXPRESSION TAG	UNP P10635
C	25	LYS	-	EXPRESSION TAG	UNP P10635
C	26	LYS	-	EXPRESSION TAG	UNP P10635
C	27	THR	-	EXPRESSION TAG	UNP P10635
C	28	SER	-	EXPRESSION TAG	UNP P10635
C	29	SER	-	EXPRESSION TAG	UNP P10635
C	30	LYS	-	EXPRESSION TAG	UNP P10635
C	31	GLY	-	EXPRESSION TAG	UNP P10635
C	32	LYS	-	EXPRESSION TAG	UNP P10635
C	33	LEU	-	EXPRESSION TAG	UNP P10635
C	498	HIS	-	EXPRESSION TAG	UNP P10635
C	499	HIS	-	EXPRESSION TAG	UNP P10635
C	500	HIS	-	EXPRESSION TAG	UNP P10635
C	501	HIS	-	EXPRESSION TAG	UNP P10635
D	23	MET	-	EXPRESSION TAG	UNP P10635
D	24	ALA	-	EXPRESSION TAG	UNP P10635
D	25	LYS	-	EXPRESSION TAG	UNP P10635
D	26	LYS	-	EXPRESSION TAG	UNP P10635
D	27	THR	-	EXPRESSION TAG	UNP P10635
D	28	SER	-	EXPRESSION TAG	UNP P10635
D	29	SER	-	EXPRESSION TAG	UNP P10635
D	30	LYS	-	EXPRESSION TAG	UNP P10635
D	31	GLY	-	EXPRESSION TAG	UNP P10635
D	32	LYS	-	EXPRESSION TAG	UNP P10635
D	33	LEU	-	EXPRESSION TAG	UNP P10635
D	498	HIS	-	EXPRESSION TAG	UNP P10635
D	499	HIS	-	EXPRESSION TAG	UNP P10635
D	500	HIS	-	EXPRESSION TAG	UNP P10635
D	501	HIS	-	EXPRESSION TAG	UNP P10635

- Molecule 2 is 10-{2-[(2R)-1-METHYLPIPERIDIN-2-YL]ETHYL}-2-(METHYLSULFANYL)-10H-PHENOTHIAZINE (three-letter code: RTZ) (formula: C₂₁H₂₆N₂S₂).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	S	0	0
			25	21	2	2		
2	A	1	Total	C	N	S	0	0
			25	21	2	2		
2	B	1	Total	C	N	S	0	0
			25	21	2	2		
2	B	1	Total	C	N	S	0	0
			25	21	2	2		
2	C	1	Total	C	N	S	0	0
			25	21	2	2		
2	C	1	Total	C	N	S	0	0
			25	21	2	2		
2	D	1	Total	C	N	S	0	0
			25	21	2	2		
2	D	1	Total	C	N	S	0	0
			25	21	2	2		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

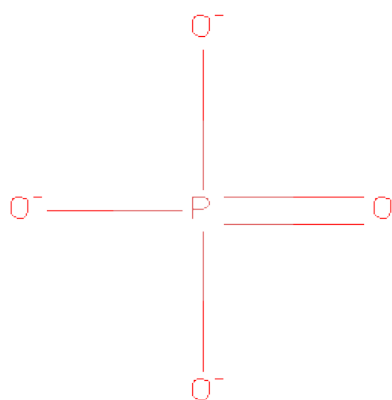
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Zn	0	0
			1	1		
3	A	3	Total	Zn	0	0
			3	3		
3	D	1	Total	Zn	0	0
			1	1		
3	C	2	Total	Zn	0	0
			2	2		

- Molecule 4 is GLYCEROL (three-letter code: GOL) (formula: $C_3H_8O_3$).



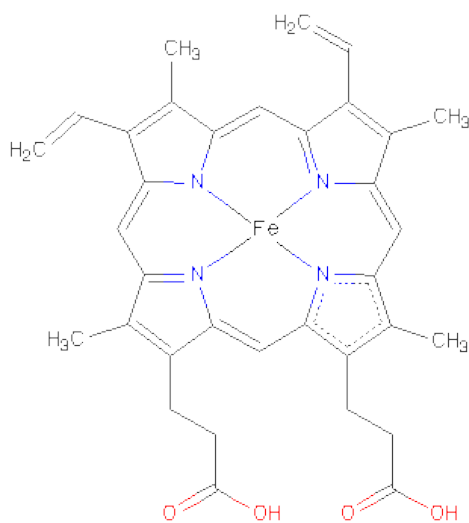
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	A	1	Total	C	O	0	0
			6	3	3		
4	B	1	Total	C	O	0	0
			6	3	3		
4	C	1	Total	C	O	0	0
			6	3	3		
4	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 5 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	O	P	0	0
			5	4	1		

- Molecule 6 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
6	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

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Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
6	C	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		
6	D	1	Total	C	Fe	N	O	0	0
			43	34	1	4	4		

- Molecule 7 is water.

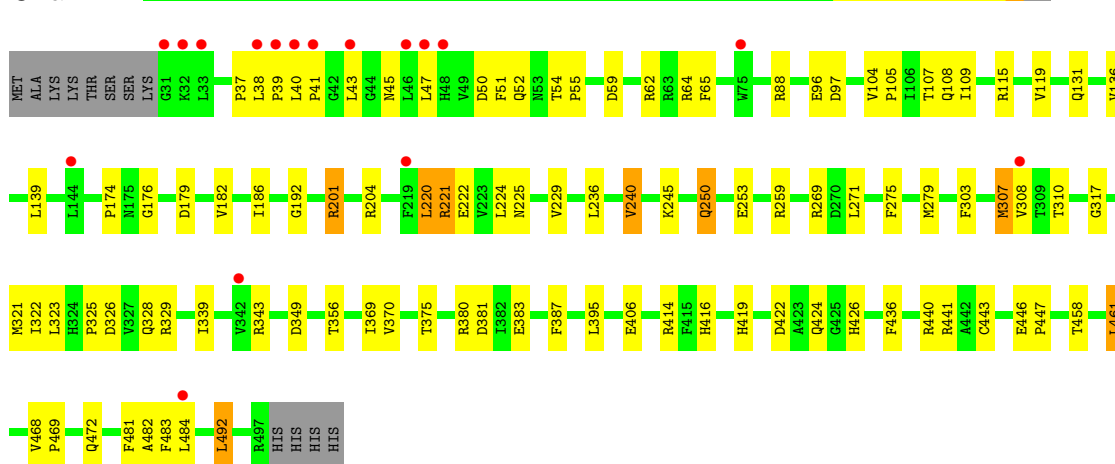
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	146	Total	O	0	0
			146	146		
7	B	109	Total	O	0	0
			109	109		
7	C	134	Total	O	0	0
			134	134		
7	D	115	Total	O	0	0
			115	115		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

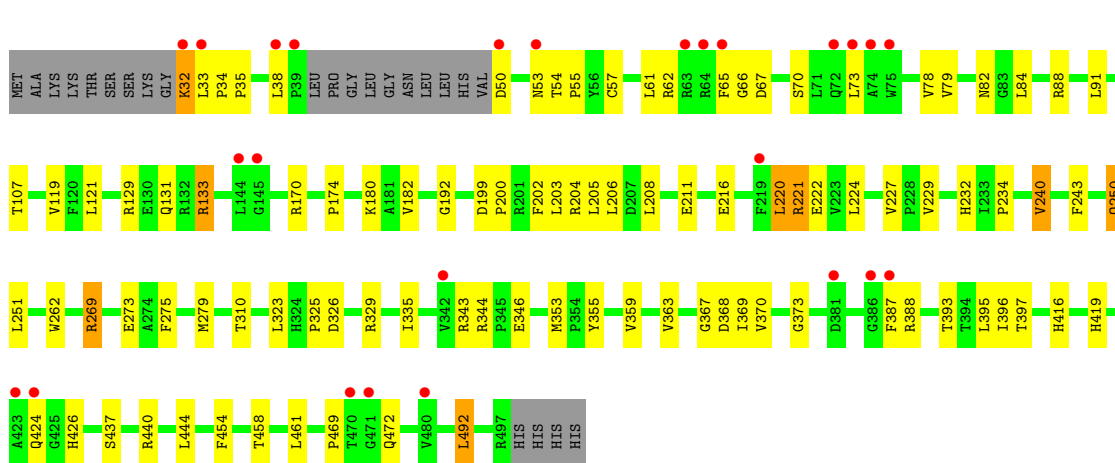
• Molecule 1: Cytochrome P450 2D6

Chain A:



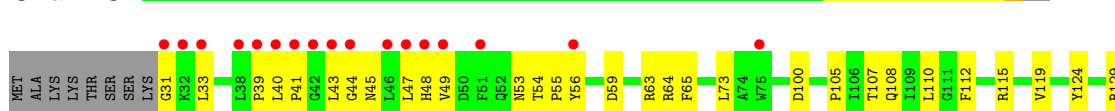
• Molecule 1: Cytochrome P450 2D6

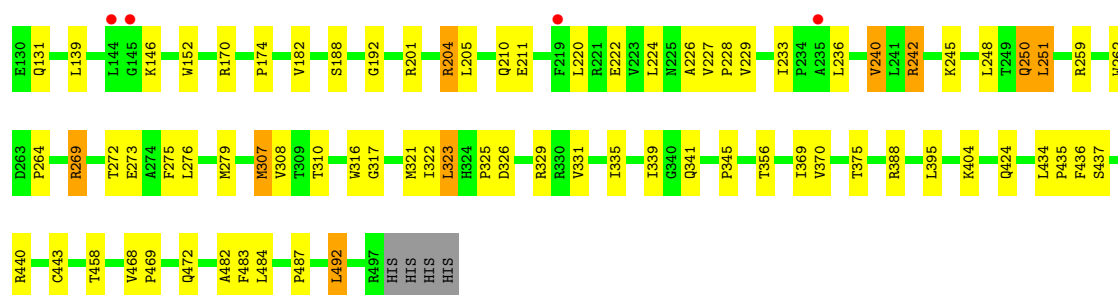
Chain B:



• Molecule 1: Cytochrome P450 2D6

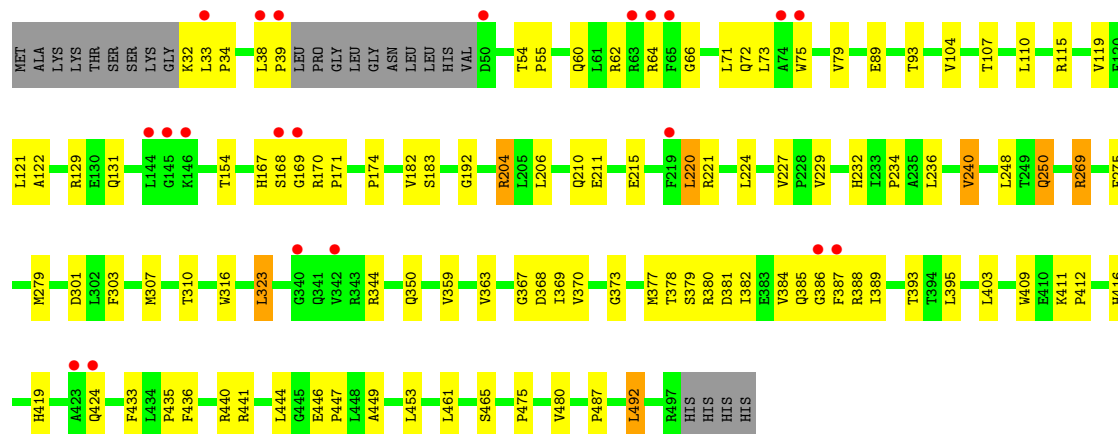
Chain C:





• Molecule 1: Cytochrome P450 2D6

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.19Å 191.96Å 249.48Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.90 – 2.10 36.90 – 2.04	Depositor EDS
% Data completeness (in resolution range)	99.5 (36.90-2.10) 99.2 (36.90-2.04)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	0.09	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.49 (at 2.05Å)	Xtriage
Refinement program	CNS 1.3	Depositor
R, R_{free}	0.222 , 0.247 0.218 , 0.244	Depositor DCC
R_{free} test set	8054 reflections (5.29%)	DCC
Wilson B-factor (Å ²)	26.9	Xtriage
Anisotropy	0.514	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 39.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	2 of 174290 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	15538	wwPDB-VP
Average B, all atoms (Å ²)	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 23.25 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 4.8395e-03. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, ZN, RTZ, PO4, HEM

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	0.36	0/3790	0.59	1/5155 (0.0%)
1	B	0.34	0/3711	0.55	0/5045
1	C	0.36	0/3790	0.57	1/5155 (0.0%)
1	D	0.34	0/3711	0.55	0/5045
All	All	0.35	0/15002	0.57	2/20400 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	441	ARG	NE-CZ-NH2	-5.93	117.34	120.30
1	C	434	LEU	CA-CB-CG	5.41	127.74	115.30

There are no chirality outliers.

There are no planarity outliers.

5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3690	0	3679	76	0
1	B	3614	0	3596	82	0
1	C	3690	0	3679	79	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	D	3614	0	3596	81	0
2	A	50	0	52	2	0
2	B	50	0	52	1	0
2	C	50	0	52	3	0
2	D	50	0	52	1	0
3	A	3	0	0	0	0
3	B	1	0	0	0	0
3	C	2	0	0	0	0
3	D	1	0	0	0	0
4	A	24	0	32	4	0
4	B	6	0	8	4	0
4	C	6	0	8	0	0
4	D	6	0	8	6	0
5	A	5	0	0	0	0
6	A	43	0	30	0	0
6	B	43	0	30	1	0
6	C	43	0	30	1	0
6	D	43	0	30	0	0
7	A	146	0	0	3	0
7	B	109	0	0	2	0
7	C	134	0	0	5	0
7	D	115	0	0	4	0
All	All	15538	0	14934	317	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 11.

All (317) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:32:LYS:HA	1:B:32:LYS:HE3	1.47	0.96
1:A:38:LEU:HD12	1:A:39:PRO:HD2	1.56	0.86
1:A:406:GLU:O	1:D:411:LYS:HE3	1.76	0.85
1:D:220:LEU:HD22	1:D:240:VAL:HG13	1.63	0.81
2:A:1:RTZ:HAI	7:A:628:HOH:O	1.79	0.81
1:B:170:ARG:HH11	1:B:170:ARG:HB2	1.49	0.78
1:B:224:LEU:HD12	1:B:240:VAL:HG11	1.66	0.78
1:A:204:ARG:HH11	1:A:204:ARG:HG2	1.49	0.77
1:D:224:LEU:HD12	1:D:240:VAL:HG11	1.66	0.77
1:D:71:LEU:HD11	1:D:73:LEU:HG	1.67	0.76
1:C:40:LEU:HG	1:C:41:PRO:HD2	1.70	0.71
1:D:221:ARG:NH2	4:D:750:GOL:O1	2.22	0.71
1:D:38:LEU:HB2	1:D:39:PRO:HD3	1.73	0.70

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:183:SER:OG	1:D:206:LEU:HD21	1.91	0.69
1:B:424:GLN:O	1:D:424:GLN:HG3	1.92	0.69
1:C:339:ILE:HD13	1:C:345:PRO:HB3	1.73	0.69
1:A:43:LEU:HD22	1:A:47:LEU:HD11	1.75	0.69
1:A:105:PRO:O	1:A:108:GLN:HG3	1.93	0.69
1:B:35:PRO:HB2	1:B:65:PHE:HD2	1.56	0.69
1:A:339:ILE:HG23	1:A:343:ARG:NH1	2.08	0.68
1:B:204:ARG:NH1	1:B:208:LEU:HD11	2.08	0.68
1:A:369:ILE:HG13	1:A:370:VAL:HG23	1.76	0.68
1:B:34:PRO:HG2	1:B:70:SER:OG	1.94	0.68
1:A:224:LEU:HD12	1:A:240:VAL:HG11	1.76	0.68
1:D:378:THR:HG23	1:D:395:LEU:HD11	1.76	0.67
1:A:307:MET:HG3	1:A:308:VAL:N	2.10	0.67
1:B:204:ARG:NH1	1:B:250:GLN:HG2	2.10	0.66
1:A:139:LEU:HD13	7:A:590:HOH:O	1.96	0.65
1:B:129:ARG:NH1	7:B:601:HOH:O	2.30	0.65
1:C:192:GLY:O	1:C:269:ARG:NH2	2.29	0.65
1:B:170:ARG:NH1	1:B:170:ARG:HB2	2.12	0.65
1:B:35:PRO:HG2	1:B:65:PHE:HB3	1.79	0.65
1:D:224:LEU:HD12	1:D:240:VAL:CG1	2.26	0.64
1:C:369:ILE:HG13	1:C:370:VAL:HG23	1.79	0.63
1:C:250:GLN:HE21	1:C:250:GLN:HA	1.62	0.63
1:C:204:ARG:HH11	1:C:204:ARG:HG2	1.63	0.63
1:C:469:PRO:HB2	1:C:472:GLN:NE2	2.14	0.62
1:A:192:GLY:O	1:A:269:ARG:NH2	2.32	0.62
1:D:204:ARG:HG2	1:D:204:ARG:HH11	1.65	0.62
1:B:133:ARG:HH11	1:B:133:ARG:HG2	1.63	0.62
1:C:222:GLU:OE1	2:C:2:RTZ:HAAA	1.99	0.62
1:C:129:ARG:NH1	7:C:4:HOH:O	2.32	0.61
1:D:104:VAL:HG11	4:D:750:GOL:H11	1.83	0.60
1:C:250:GLN:CA	1:C:250:GLN:HE21	2.14	0.60
1:B:227:VAL:HG12	1:B:229:VAL:HG12	1.82	0.60
1:B:202:PHE:O	1:B:206:LEU:HD13	2.02	0.60
1:D:174:PRO:HD3	1:D:492:LEU:HD22	1.84	0.60
1:C:307:MET:HG3	1:C:308:VAL:N	2.16	0.59
1:D:227:VAL:HG12	1:D:229:VAL:HG12	1.83	0.59
1:C:224:LEU:HD12	1:C:240:VAL:HG11	1.84	0.58
1:D:369:ILE:HG13	1:D:370:VAL:HG23	1.84	0.58
1:D:54:THR:HB	1:D:55:PRO:HD3	1.83	0.58
1:C:39:PRO:HA	1:C:45:ASN:ND2	2.17	0.58
1:A:224:LEU:HD12	1:A:240:VAL:CG1	2.32	0.58
1:D:359:VAL:O	1:D:363:VAL:HG23	2.04	0.57

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:204:ARG:HG2	1:A:204:ARG:NH1	2.18	0.57
1:D:38:LEU:HB2	1:D:39:PRO:CD	2.35	0.57
1:D:170:ARG:NH1	1:D:170:ARG:HB2	2.20	0.56
1:C:233:ILE:HD12	1:C:233:ILE:N	2.19	0.56
1:C:139:LEU:HD13	7:C:623:HOH:O	2.05	0.56
1:C:224:LEU:HD12	1:C:240:VAL:CG1	2.35	0.56
1:B:204:ARG:HH11	1:B:250:GLN:HG2	1.70	0.56
1:B:182:VAL:HG11	1:B:310:THR:HB	1.87	0.56
1:D:236:LEU:O	1:D:240:VAL:HG22	2.06	0.56
1:B:174:PRO:HD3	1:B:492:LEU:HD22	1.86	0.56
1:A:422:ASP:HB3	1:C:424:GLN:HG3	1.87	0.56
1:A:220:LEU:HD13	1:A:240:VAL:HG13	1.87	0.56
1:D:370:VAL:HG12	1:D:373:GLY:HA2	1.88	0.56
1:A:136:VAL:HG13	4:A:753:GOL:H12	1.88	0.56
1:D:33:LEU:HD11	1:D:388:ARG:NH1	2.21	0.56
1:C:182:VAL:HG11	1:C:310:THR:HB	1.88	0.56
1:A:109:ILE:CG2	1:A:245:LYS:HD2	2.36	0.55
1:B:355:TYR:O	1:B:359:VAL:HG23	2.07	0.55
1:D:232:HIS:O	1:D:234:PRO:HD3	2.07	0.55
1:B:54:THR:HB	1:B:55:PRO:HD3	1.88	0.55
1:B:84:LEU:HD11	1:B:88:ARG:HD2	1.88	0.55
1:B:369:ILE:HG13	1:B:370:VAL:HG23	1.88	0.54
1:D:250:GLN:HA	1:D:250:GLN:HE21	1.72	0.54
1:B:192:GLY:O	1:B:269:ARG:NH2	2.41	0.54
1:C:227:VAL:HG12	1:C:229:VAL:HG12	1.88	0.54
1:B:91:LEU:HD11	1:B:397:THR:HG21	1.88	0.54
1:C:40:LEU:CD2	1:C:47:LEU:HD12	2.38	0.54
1:A:469:PRO:HB2	1:A:472:GLN:NE2	2.22	0.54
1:B:38:LEU:N	1:B:38:LEU:HD22	2.23	0.54
1:A:482:ALA:O	1:A:483:PHE:HB3	2.07	0.54
1:A:96:GLU:HG2	1:A:440:ARG:NH1	2.22	0.54
1:B:221:ARG:HG3	1:B:222:GLU:N	2.23	0.54
1:B:121:LEU:CD1	4:B:750:GOL:H11	2.38	0.54
1:A:174:PRO:HD3	1:A:492:LEU:HD22	1.90	0.54
1:A:43:LEU:N	1:A:43:LEU:HD12	2.24	0.53
1:B:32:LYS:HE3	1:B:32:LYS:CA	2.31	0.53
1:A:40:LEU:HG	1:A:41:PRO:HD2	1.89	0.53
1:A:97:ASP:OD2	1:A:380:ARG:HD2	2.09	0.53
1:D:384:VAL:HG12	1:D:385:GLN:HG3	1.90	0.53
1:C:119:VAL:HG22	1:C:131:GLN:HB3	1.90	0.53
1:A:236:LEU:O	1:A:240:VAL:HG22	2.08	0.53
1:B:262:TRP:CD1	1:B:273:GLU:HG2	2.44	0.53

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:50:ASP:OD1	1:A:52:GLN:HG2	2.07	0.53
1:D:60:GLN:O	1:D:64:ARG:HG3	2.08	0.53
1:B:454:PHE:O	1:B:458:THR:HG23	2.09	0.53
1:A:322:ILE:HD12	1:A:468:VAL:HG12	1.91	0.53
1:B:199:ASP:O	1:B:203:LEU:HG	2.09	0.52
1:A:104:VAL:HG11	4:A:750:GOL:H11	1.92	0.52
1:A:326:ASP:N	1:A:326:ASP:OD2	2.42	0.52
1:C:31:GLY:HA2	1:C:388:ARG:HD3	1.92	0.52
1:D:379:SER:O	1:D:380:ARG:HG3	2.10	0.52
1:C:440:ARG:HH21	1:C:440:ARG:HG3	1.74	0.52
1:C:40:LEU:HB2	1:C:44:GLY:O	2.09	0.52
1:D:170:ARG:HH11	1:D:170:ARG:HB2	1.73	0.52
1:B:204:ARG:HH11	1:B:250:GLN:CG	2.22	0.52
1:B:62:ARG:HD2	1:B:82:ASN:HB3	1.91	0.52
1:D:440:ARG:HG3	7:D:609:HOH:O	2.10	0.52
1:B:62:ARG:HH11	1:B:82:ASN:HB2	1.75	0.51
1:D:275:PHE:CE2	1:D:279:MET:HE2	2.45	0.51
1:B:359:VAL:O	1:B:363:VAL:HG23	2.10	0.51
1:D:119:VAL:CG2	1:D:131:GLN:HB3	2.40	0.51
1:C:110:LEU:HD22	1:C:248:LEU:HD12	1.91	0.51
1:A:37:PRO:HG2	1:A:45:ASN:HD22	1.76	0.51
1:D:71:LEU:C	1:D:71:LEU:HD12	2.31	0.51
1:C:49:VAL:HG21	1:C:73:LEU:HD21	1.93	0.51
1:D:182:VAL:HG11	1:D:310:THR:HB	1.92	0.51
1:A:339:ILE:HG23	1:A:343:ARG:HH12	1.75	0.51
1:B:440:ARG:HG3	7:B:548:HOH:O	2.10	0.51
1:C:325:PRO:O	1:C:329:ARG:HG3	2.12	0.50
1:B:224:LEU:HD12	1:B:240:VAL:CG1	2.39	0.50
1:A:422:ASP:CB	1:C:424:GLN:HG3	2.41	0.50
1:C:469:PRO:O	1:C:472:GLN:HB2	2.12	0.50
1:B:170:ARG:HH11	1:B:170:ARG:CB	2.24	0.50
1:A:221:ARG:HG3	1:A:222:GLU:N	2.27	0.50
1:C:119:VAL:CG2	1:C:131:GLN:HB3	2.42	0.49
1:A:186:ILE:HD13	1:A:303:PHE:HA	1.94	0.49
1:C:40:LEU:CG	1:C:41:PRO:HD2	2.42	0.49
1:D:221:ARG:HH22	4:D:750:GOL:C1	2.25	0.49
1:A:105:PRO:HD2	1:A:225:ASN:OD1	2.12	0.49
1:D:129:ARG:NH2	7:D:512:HOH:O	2.45	0.49
1:C:45:ASN:HA	1:C:48:HIS:ND1	2.27	0.49
1:A:328:GLN:HG3	1:A:461:LEU:CD1	2.42	0.49
1:B:121:LEU:HD11	4:B:750:GOL:H11	1.95	0.49
1:A:325:PRO:O	1:A:329:ARG:HG3	2.12	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:416:HIS:HB3	1:A:419:HIS:CE1	2.48	0.49
1:B:275:PHE:CE2	1:B:279:MET:HE2	2.48	0.49
1:A:59:ASP:HA	1:A:62:ARG:NH1	2.27	0.49
1:D:89:GLU:O	1:D:93:THR:HB	2.12	0.48
1:A:250:GLN:HE21	1:A:250:GLN:CA	2.25	0.48
1:C:54:THR:HB	1:C:55:PRO:HD3	1.94	0.48
1:A:424:GLN:HG2	1:A:426:HIS:CE1	2.49	0.48
1:B:335:ILE:HG12	1:B:353:MET:HE1	1.94	0.48
1:C:236:LEU:O	1:C:240:VAL:HG22	2.13	0.48
1:A:64:ARG:HD2	1:A:65:PHE:CE1	2.48	0.48
1:C:210:GLN:NE2	1:C:211:GLU:HG2	2.28	0.48
1:D:121:LEU:CD1	4:D:750:GOL:H12	2.44	0.48
1:A:356:THR:HG21	1:A:458:THR:HG22	1.96	0.48
1:A:40:LEU:HD23	1:A:41:PRO:N	2.29	0.48
1:D:168:SER:C	1:D:170:ARG:H	2.17	0.48
1:B:88:ARG:HB3	1:B:88:ARG:NH1	2.29	0.48
1:A:119:VAL:CG2	1:A:131:GLN:HB3	2.43	0.48
1:A:229:VAL:HG13	7:A:604:HOH:O	2.13	0.48
1:A:414:ARG:HD3	7:D:595:HOH:O	2.13	0.48
1:C:356:THR:HG21	1:C:458:THR:HG22	1.94	0.47
1:C:482:ALA:O	1:C:483:PHE:CB	2.62	0.47
1:D:122:ALA:O	1:D:441:ARG:NH2	2.37	0.47
1:A:107:THR:HG22	4:A:750:GOL:H32	1.96	0.47
1:C:210:GLN:HB3	7:C:603:HOH:O	2.15	0.47
1:A:275:PHE:CE2	1:A:279:MET:HE2	2.49	0.47
1:B:78:VAL:HG11	1:B:396:ILE:HD12	1.96	0.47
1:A:37:PRO:HG2	1:A:45:ASN:ND2	2.29	0.47
1:B:275:PHE:HE2	1:B:279:MET:HE2	1.79	0.47
1:A:326:ASP:HA	1:A:329:ARG:NH2	2.30	0.47
1:D:210:GLN:HG3	1:D:211:GLU:N	2.29	0.47
1:C:40:LEU:HD22	1:C:47:LEU:HD12	1.97	0.47
1:B:326:ASP:N	1:B:326:ASP:OD2	2.47	0.47
1:B:326:ASP:HA	1:B:329:ARG:NH1	2.30	0.46
1:D:110:LEU:HD22	1:D:248:LEU:HD12	1.97	0.46
1:D:220:LEU:HD22	1:D:240:VAL:CG1	2.39	0.46
1:A:40:LEU:HD22	1:A:43:LEU:HD13	1.96	0.46
1:A:383:GLU:HA	1:A:387:PHE:O	2.15	0.46
1:B:119:VAL:HG22	1:B:131:GLN:HB3	1.97	0.46
1:D:119:VAL:HG22	1:D:131:GLN:HB3	1.96	0.46
1:D:71:LEU:CD1	1:D:73:LEU:HG	2.39	0.46
1:A:88:ARG:NH1	1:A:88:ARG:HB3	2.31	0.46
1:B:416:HIS:HB3	1:B:419:HIS:CE1	2.51	0.46

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:307:MET:CG	1:A:308:VAL:N	2.78	0.46
1:C:262:TRP:CD1	1:C:273:GLU:HG2	2.50	0.46
1:C:174:PRO:HD3	1:C:492:LEU:HD22	1.98	0.46
1:D:435:PRO:HG2	1:D:436:PHE:CD1	2.50	0.46
1:D:107:THR:HG22	4:D:750:GOL:H32	1.96	0.46
1:C:204:ARG:NH1	1:C:204:ARG:HG2	2.28	0.46
1:C:436:PHE:HB3	1:C:443:CYS:HB3	1.97	0.46
1:D:192:GLY:O	1:D:269:ARG:NH2	2.48	0.46
1:B:211:GLU:HB3	1:B:243:PHE:CD2	2.51	0.46
1:B:35:PRO:HD3	1:B:387:PHE:CE1	2.51	0.46
1:C:205:LEU:HD11	1:C:251:LEU:HD13	1.98	0.46
1:C:245:LYS:HE2	1:C:245:LYS:HB3	1.84	0.45
1:B:33:LEU:HD23	1:B:388:ARG:O	2.16	0.45
1:B:444:LEU:C	1:B:444:LEU:HD12	2.37	0.45
1:C:275:PHE:CE2	1:C:279:MET:HE2	2.51	0.45
1:B:220:LEU:HD22	1:B:240:VAL:HG13	1.99	0.45
1:C:316:TRP:CD1	1:C:487:PRO:HD3	2.51	0.45
1:C:375:THR:HG23	2:C:2:RTZ:CAJ	2.47	0.45
1:D:370:VAL:HG12	1:D:373:GLY:CA	2.47	0.45
1:D:121:LEU:HD11	4:D:750:GOL:H12	1.98	0.45
1:D:275:PHE:HE2	1:D:279:MET:HE2	1.81	0.45
1:D:367:GLY:O	1:D:368:ASP:C	2.56	0.45
1:D:368:ASP:OD1	1:D:403:LEU:HD12	2.16	0.45
1:A:375:THR:HG23	2:A:2:RTZ:CAJ	2.47	0.45
1:C:226:ALA:C	1:C:228:PRO:HD3	2.37	0.44
1:A:303:PHE:O	1:A:307:MET:HB3	2.18	0.44
1:D:62:ARG:HA	1:D:66:GLY:O	2.17	0.44
1:C:33:LEU:HD23	1:C:388:ARG:O	2.18	0.44
1:D:168:SER:O	1:D:170:ARG:N	2.51	0.44
1:D:433:PHE:CZ	1:D:435:PRO:HG3	2.52	0.44
1:A:54:THR:HB	1:A:55:PRO:HD3	2.00	0.44
1:A:271:LEU:C	1:A:271:LEU:HD23	2.38	0.44
1:D:71:LEU:HD12	1:D:72:GLN:N	2.33	0.44
1:A:179:ASP:HB3	1:A:307:MET:HE1	1.99	0.44
1:D:416:HIS:HB3	1:D:419:HIS:CE1	2.53	0.44
1:C:100:ASP:HA	1:C:124:TYR:HB2	2.00	0.44
1:D:382:ILE:HG13	1:D:389:ILE:HB	1.99	0.44
1:C:251:LEU:HD12	1:C:251:LEU:HA	1.87	0.44
1:A:436:PHE:HB3	1:A:443:CYS:HB3	1.98	0.44
1:C:404:LYS:NZ	7:C:586:HOH:O	2.49	0.43
1:B:424:GLN:CD	1:B:424:GLN:H	2.22	0.43
1:B:180:LYS:HE2	1:B:203:LEU:CD2	2.48	0.43

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:C:492:LEU:O	1:C:492:LEU:HD23	2.18	0.43
1:D:75:TRP:N	1:D:75:TRP:CD1	2.86	0.43
1:D:377:MET:HA	1:D:393:THR:O	2.18	0.43
1:B:220:LEU:HD13	1:B:240:VAL:HG13	2.01	0.43
1:D:204:ARG:NH1	1:D:204:ARG:HG2	2.32	0.43
1:A:201:ARG:HD2	1:A:253:GLU:OE2	2.18	0.43
1:C:119:VAL:HG22	1:C:131:GLN:CB	2.47	0.43
1:B:62:ARG:HA	1:B:66:GLY:O	2.18	0.43
1:B:205:LEU:HD11	1:B:251:LEU:HD13	2.01	0.43
1:C:250:GLN:NE2	1:C:250:GLN:HA	2.32	0.43
1:D:79:VAL:HG21	1:D:389:ILE:HD12	2.01	0.43
1:C:331:VAL:O	1:C:335:ILE:HG13	2.18	0.43
1:D:446:GLU:HB3	1:D:447:PRO:HD3	2.00	0.43
1:D:444:LEU:HD12	1:D:444:LEU:C	2.39	0.43
1:A:119:VAL:HG22	1:A:131:GLN:HB3	2.00	0.43
1:D:154:THR:HG22	1:D:344:ARG:HD3	2.01	0.43
1:C:53:ASN:HB3	1:C:56:TYR:HD2	1.84	0.43
1:B:78:VAL:HG11	1:B:396:ILE:CD1	2.49	0.43
1:A:51:PHE:HA	1:A:54:THR:OG1	2.19	0.43
1:B:437:SER:HB3	6:B:800:HEM:HBA1	2.00	0.43
1:A:317:GLY:O	1:A:321:MET:HG2	2.19	0.43
1:B:269:ARG:HG2	1:B:273:GLU:OE1	2.19	0.43
1:D:350:GLN:NE2	7:D:603:HOH:O	2.52	0.43
1:B:232:HIS:O	1:B:234:PRO:HD3	2.18	0.43
1:A:469:PRO:O	1:A:472:GLN:HB2	2.19	0.42
1:D:167:HIS:O	1:D:170:ARG:HB3	2.20	0.42
1:D:33:LEU:HD12	1:D:33:LEU:N	2.34	0.42
1:C:40:LEU:HD13	1:C:48:HIS:HE1	1.85	0.42
1:A:176:GLY:HA2	4:A:751:GOL:H11	2.00	0.42
1:B:424:GLN:HB2	1:B:426:HIS:ND1	2.34	0.42
1:C:152:TRP:CD2	1:C:188:SER:HB3	2.55	0.42
1:D:449:ALA:O	1:D:453:LEU:HG	2.19	0.42
1:D:303:PHE:O	1:D:307:MET:HB2	2.19	0.42
1:A:37:PRO:O	1:A:45:ASN:ND2	2.51	0.42
1:A:182:VAL:HG11	1:A:310:THR:HB	2.02	0.42
1:C:323:LEU:HA	1:C:323:LEU:HD12	1.85	0.42
1:B:107:THR:HG22	4:B:750:GOL:H32	2.01	0.42
1:B:62:ARG:NH1	1:B:67:ASP:OD2	2.52	0.42
1:C:226:ALA:O	1:C:228:PRO:HD3	2.19	0.42
1:C:262:TRP:O	1:C:264:PRO:HD3	2.20	0.42
1:B:57:CYS:O	1:B:61:LEU:HG	2.20	0.42
1:C:326:ASP:OD2	1:C:326:ASP:N	2.52	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:32:LYS:HD2	1:D:386:GLY:O	2.20	0.42
1:D:301:ASP:OD2	2:D:1:RTZ:NAV	2.53	0.42
1:B:54:THR:N	1:B:55:PRO:CD	2.82	0.42
1:B:119:VAL:CG2	1:B:131:GLN:HB3	2.50	0.42
1:C:435:PRO:HG2	1:C:436:PHE:CD1	2.55	0.42
1:C:64:ARG:HD2	1:C:65:PHE:CE1	2.55	0.42
1:B:343:ARG:HH21	1:B:346:GLU:CD	2.23	0.41
1:C:105:PRO:O	1:C:108:GLN:HG3	2.20	0.41
1:B:133:ARG:NH1	1:B:133:ARG:HG2	2.33	0.41
1:A:446:GLU:HB3	1:A:447:PRO:HD3	2.02	0.41
1:D:215:GLU:O	1:D:221:ARG:NH1	2.52	0.41
1:B:204:ARG:NH1	1:B:250:GLN:CG	2.81	0.41
1:A:40:LEU:CG	1:A:41:PRO:HD2	2.50	0.41
1:D:409:TRP:O	1:D:412:PRO:HD3	2.21	0.41
1:C:272:THR:O	1:C:276:LEU:HG	2.20	0.41
1:C:322:ILE:HD12	1:C:468:VAL:HG12	2.02	0.41
1:A:481:PHE:CZ	1:A:483:PHE:HA	2.55	0.41
1:C:482:ALA:O	1:C:483:PHE:HB3	2.21	0.41
2:B:1:RTZ:HAG	2:B:1:RTZ:HAB	2.02	0.41
1:C:43:LEU:CD1	1:C:47:LEU:HD11	2.51	0.41
1:B:370:VAL:HG12	1:B:373:GLY:HA2	2.01	0.41
1:C:107:THR:HB	1:C:112:PHE:CG	2.56	0.41
1:B:62:ARG:HH11	1:B:62:ARG:HG3	1.85	0.41
1:B:325:PRO:O	1:B:329:ARG:HG3	2.20	0.41
1:B:367:GLY:O	1:B:368:ASP:C	2.57	0.41
1:A:343:ARG:NH2	1:A:349:ASP:OD2	2.54	0.41
1:C:307:MET:CG	1:C:308:VAL:N	2.82	0.41
1:B:216:GLU:OE2	4:B:750:GOL:H12	2.20	0.41
1:B:50:ASP:OD2	1:B:53:ASN:ND2	2.49	0.41
1:D:316:TRP:CD1	1:D:487:PRO:HD3	2.55	0.41
1:B:469:PRO:HB2	1:B:472:GLN:CD	2.41	0.41
1:D:171:PRO:HA	1:D:492:LEU:O	2.21	0.41
1:C:317:GLY:O	1:C:321:MET:HG2	2.21	0.41
1:C:242:ARG:HB2	1:C:242:ARG:HH11	1.86	0.41
1:B:79:VAL:HG23	1:B:393:THR:HG21	2.03	0.40
1:C:437:SER:HB3	6:C:800:HEM:HBA1	2.02	0.40
1:D:480:VAL:HG13	1:D:480:VAL:O	2.22	0.40
1:D:34:PRO:HG3	1:D:387:PHE:HB3	2.03	0.40
1:A:307:MET:HG3	1:A:308:VAL:H	1.84	0.40
1:B:199:ASP:HA	1:B:200:PRO:HD3	1.96	0.40
1:B:73:LEU:HD12	1:B:78:VAL:HG21	2.02	0.40
1:D:323:LEU:CD1	1:D:475:PRO:HG2	2.51	0.40

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:71:LEU:HD11	1:D:73:LEU:CG	2.42	0.40
1:A:492:LEU:O	1:A:492:LEU:HD23	2.22	0.40
1:C:468:VAL:O	1:C:468:VAL:HG23	2.21	0.40
1:C:341:GLN:HG2	1:C:341:GLN:O	2.21	0.40
1:B:461:LEU:HD23	1:B:461:LEU:HA	1.84	0.40
1:A:40:LEU:HD22	1:A:43:LEU:H	1.87	0.40
2:C:2:RTZ:HAA	7:C:613:HOH:O	2.21	0.40
1:C:440:ARG:NH2	1:C:440:ARG:HG3	2.37	0.40
1:B:62:ARG:NH1	1:B:82:ASN:HB2	2.36	0.40
1:D:461:LEU:HA	1:D:461:LEU:HD23	1.92	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	465/479 (97%)	452 (97%)	13 (3%)	0	100	100
1	B	452/479 (94%)	437 (97%)	15 (3%)	0	100	100
1	C	465/479 (97%)	449 (97%)	16 (3%)	0	100	100
1	D	452/479 (94%)	435 (96%)	16 (4%)	1 (0%)	56	57
All	All	1834/1916 (96%)	1773 (97%)	60 (3%)	1 (0%)	59	61

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	169	GLY

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	398/409 (97%)	384 (96%)	14 (4%)	48	48
1	B	390/409 (95%)	379 (97%)	11 (3%)	56	59
1	C	398/409 (97%)	379 (95%)	19 (5%)	35	32
1	D	390/409 (95%)	380 (97%)	10 (3%)	59	62
All	All	1576/1636 (96%)	1522 (97%)	54 (3%)	49	49

All (54) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	115	ARG
1	A	201	ARG
1	A	220	LEU
1	A	221	ARG
1	A	240	VAL
1	A	250	GLN
1	A	259	ARG
1	A	307	MET
1	A	323	LEU
1	A	381	ASP
1	A	395	LEU
1	A	461	LEU
1	A	484	LEU
1	A	492	LEU
1	B	32	LYS
1	B	133	ARG
1	B	220	LEU
1	B	221	ARG
1	B	240	VAL
1	B	250	GLN
1	B	269	ARG
1	B	323	LEU
1	B	344	ARG
1	B	395	LEU
1	B	492	LEU
1	C	59	ASP
1	C	63	ARG
1	C	115	ARG
1	C	146	LYS
1	C	170	ARG
1	C	201	ARG

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Mol	Chain	Res	Type
1	C	204	ARG
1	C	220	LEU
1	C	240	VAL
1	C	242	ARG
1	C	250	GLN
1	C	251	LEU
1	C	259	ARG
1	C	269	ARG
1	C	307	MET
1	C	323	LEU
1	C	395	LEU
1	C	484	LEU
1	C	492	LEU
1	D	115	ARG
1	D	204	ARG
1	D	220	LEU
1	D	240	VAL
1	D	250	GLN
1	D	269	ARG
1	D	323	LEU
1	D	381	ASP
1	D	465	SER
1	D	492	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	463	HIS
1	B	472	GLN
1	C	250	GLN
1	C	472	GLN
1	D	60	GLN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

Of 27 ligands modelled in this entry, 7 are monoatomic - leaving 20 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$
2	RTZ	A	1	-	28,28,28	1.80	7 (25%)	39,39,39	1.54	5 (12%)
2	RTZ	A	2	-	28,28,28	1.92	9 (32%)	39,39,39	1.66	3 (7%)
4	GOL	A	750	-	5,5,5	0.25	0	5,5,5	0.33	0
4	GOL	A	751	-	5,5,5	0.23	0	5,5,5	0.25	0
4	GOL	A	752	-	5,5,5	0.23	0	5,5,5	0.24	0
4	GOL	A	753	-	5,5,5	0.30	0	5,5,5	0.28	0
5	PO4	A	790	-	4,4,4	0.86	0	6,6,6	0.31	0
6	HEM	A	800	1,7	49,50,50	4.17	21 (42%)	46,82,82	1.67	9 (19%)
2	RTZ	B	1	-	28,28,28	1.79	9 (32%)	39,39,39	1.44	5 (12%)
2	RTZ	B	2	-	28,28,28	1.78	9 (32%)	39,39,39	1.18	4 (10%)
4	GOL	B	750	-	5,5,5	0.20	0	5,5,5	0.30	0
6	HEM	B	800	1,7	49,50,50	4.32	22 (44%)	46,82,82	1.73	10 (21%)
2	RTZ	C	1	-	28,28,28	1.71	8 (28%)	39,39,39	1.52	5 (12%)
2	RTZ	C	2	-	28,28,28	1.94	10 (35%)	39,39,39	1.72	4 (10%)
4	GOL	C	750	-	5,5,5	0.26	0	5,5,5	0.40	0
6	HEM	C	800	1,7	49,50,50	3.80	21 (42%)	46,82,82	1.63	10 (21%)
2	RTZ	D	1	-	28,28,28	1.82	7 (25%)	39,39,39	1.42	5 (12%)
2	RTZ	D	2	-	28,28,28	1.76	9 (32%)	39,39,39	1.18	5 (12%)
4	GOL	D	750	-	5,5,5	0.22	0	5,5,5	0.45	0
6	HEM	D	800	1	49,50,50	4.15	22 (44%)	46,82,82	1.67	8 (17%)

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
2	RTZ	A	1	-	-	0/7/30/30	0/1/4/4
2	RTZ	A	2	-	-	0/7/30/30	0/1/4/4
4	GOL	A	750	-	-	0/4/4/4	0/0/0/0
4	GOL	A	751	-	-	0/4/4/4	0/0/0/0
4	GOL	A	752	-	-	0/4/4/4	0/0/0/0
4	GOL	A	753	-	-	0/4/4/4	0/0/0/0
5	PO4	A	790	-	-	0/0/0/0	0/0/0/0
6	HEM	A	800	1,7	-	0/14/114/114	0/0/8/8
2	RTZ	B	1	-	-	0/7/30/30	0/1/4/4
2	RTZ	B	2	-	-	0/7/30/30	0/1/4/4
4	GOL	B	750	-	-	0/4/4/4	0/0/0/0
6	HEM	B	800	1,7	-	0/14/114/114	0/0/8/8
2	RTZ	C	1	-	-	0/7/30/30	0/1/4/4
2	RTZ	C	2	-	-	0/7/30/30	0/1/4/4
4	GOL	C	750	-	-	0/4/4/4	0/0/0/0
6	HEM	C	800	1,7	-	0/14/114/114	0/0/8/8
2	RTZ	D	1	-	-	0/7/30/30	0/1/4/4
2	RTZ	D	2	-	-	0/7/30/30	0/1/4/4
4	GOL	D	750	-	-	0/4/4/4	0/0/0/0
6	HEM	D	800	1	-	0/14/114/114	0/0/8/8

All (154) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	A	800	HEM	C3D-C4D	-19.61	1.39	1.44
6	B	800	HEM	C2D-C1D	-18.08	1.40	1.44
6	D	800	HEM	C2D-C1D	-17.96	1.40	1.44
6	B	800	HEM	C3D-C4D	-16.46	1.40	1.44
6	C	800	HEM	C3D-C4D	-16.10	1.40	1.44
6	D	800	HEM	C3D-C4D	-15.65	1.40	1.44
6	C	800	HEM	C2D-C1D	-15.15	1.40	1.44
6	A	800	HEM	C2D-C1D	-14.41	1.41	1.44
6	B	800	HEM	C2B-C1B	-10.97	1.41	1.44
6	D	800	HEM	C2B-C1B	-8.96	1.42	1.44
6	A	800	HEM	C2B-C1B	-7.55	1.42	1.44
6	C	800	HEM	C2B-C1B	-4.48	1.43	1.44
6	A	800	HEM	C1C-NC	-4.14	1.32	1.38
6	C	800	HEM	C3D-C2D	-4.10	1.36	1.43
6	B	800	HEM	C1C-NC	-4.07	1.32	1.38
6	A	800	HEM	C3D-C2D	-3.91	1.36	1.43
6	B	800	HEM	C3D-C2D	-3.89	1.37	1.43
6	B	800	HEM	CMA-C3A	3.89	1.59	1.51

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	800	HEM	C1C-NC	-3.84	1.32	1.38
6	D	800	HEM	C3D-C2D	-3.83	1.37	1.43
6	D	800	HEM	C1C-NC	-3.70	1.32	1.38
6	A	800	HEM	C1D-ND	-3.56	1.29	1.37
6	D	800	HEM	FE-NA	3.48	2.07	1.92
6	B	800	HEM	FE-NA	3.48	2.07	1.92
6	D	800	HEM	CMA-C3A	3.47	1.58	1.51
6	C	800	HEM	CMC-C2C	3.46	1.58	1.47
2	C	2	RTZ	CAR-CAT	3.41	1.44	1.40
2	A	2	RTZ	CAT-SAY	3.40	1.82	1.76
6	C	800	HEM	CMA-C3A	3.40	1.58	1.51
6	D	800	HEM	CBB-CAB	3.39	1.48	1.28
6	A	800	HEM	CMA-C3A	3.39	1.58	1.51
6	B	800	HEM	FE-ND	3.38	2.10	1.97
6	C	800	HEM	CBB-CAB	3.36	1.48	1.28
6	C	800	HEM	FE-NA	3.31	2.06	1.92
6	A	800	HEM	FE-NA	3.31	2.06	1.92
6	A	800	HEM	CBB-CAB	3.30	1.48	1.28
6	B	800	HEM	CBB-CAB	3.29	1.48	1.28
2	B	2	RTZ	CAR-CAT	3.29	1.44	1.40
2	C	2	RTZ	CAT-SAY	3.28	1.82	1.76
6	A	800	HEM	CMC-C2C	3.27	1.57	1.47
2	A	2	RTZ	CAR-CAT	3.26	1.44	1.40
2	D	2	RTZ	CAR-CAT	3.26	1.44	1.40
6	D	800	HEM	CMC-C2C	3.22	1.57	1.47
6	C	800	HEM	FE-ND	3.19	2.09	1.97
2	A	1	RTZ	CAS-CAU	3.18	1.44	1.40
2	A	1	RTZ	CAR-CAT	3.17	1.44	1.40
2	D	1	RTZ	CAT-SAY	3.16	1.82	1.76
2	D	1	RTZ	CAS-CAU	3.16	1.44	1.40
6	B	800	HEM	C1B-NB	-3.15	1.33	1.39
2	D	1	RTZ	CAR-CAT	3.14	1.44	1.40
2	B	1	RTZ	CAS-CAU	3.13	1.44	1.40
6	A	800	HEM	FE-ND	3.13	2.09	1.97
6	D	800	HEM	C1B-NB	-3.11	1.33	1.39
6	C	800	HEM	C1D-ND	-3.09	1.30	1.37
2	D	2	RTZ	CAT-SAY	3.10	1.82	1.76
6	A	800	HEM	C1B-NB	-3.10	1.33	1.39
2	C	2	RTZ	CAS-CAU	3.08	1.44	1.40
6	C	800	HEM	C1B-NB	-3.07	1.33	1.39
6	D	800	HEM	CHB-C1B	3.06	1.40	1.35
6	D	800	HEM	C1D-ND	-3.06	1.30	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1	RTZ	CAR-CAT	3.06	1.44	1.40
2	A	2	RTZ	CAS-CAU	3.04	1.44	1.40
6	B	800	HEM	C1D-ND	-3.03	1.30	1.37
6	B	800	HEM	CMC-C2C	3.01	1.56	1.47
2	A	1	RTZ	CAT-SAY	3.00	1.82	1.76
2	C	1	RTZ	CAS-CAU	2.95	1.44	1.40
6	D	800	HEM	FE-ND	2.95	2.08	1.97
2	B	1	RTZ	CAT-SAY	2.95	1.82	1.76
6	D	800	HEM	C4D-ND	-2.94	1.33	1.39
2	B	2	RTZ	CAM-NAV	2.94	1.53	1.47
6	A	800	HEM	CMD-C2D	2.92	1.56	1.47
2	D	2	RTZ	CAS-CAU	2.87	1.44	1.40
2	C	1	RTZ	CAR-CAT	2.86	1.44	1.40
2	C	2	RTZ	CAL-CAN	2.86	1.57	1.52
6	B	800	HEM	C4D-ND	-2.85	1.33	1.39
2	D	2	RTZ	CAM-NAV	2.81	1.52	1.47
6	A	800	HEM	FE-NB	2.81	2.08	1.97
6	C	800	HEM	CHB-C1B	2.78	1.39	1.35
2	C	1	RTZ	CAT-SAY	2.77	1.81	1.76
2	B	2	RTZ	CAT-SAY	2.77	1.81	1.76
2	C	1	RTZ	CAM-NAV	2.76	1.52	1.47
2	B	2	RTZ	CAS-CAU	2.75	1.43	1.40
2	A	1	RTZ	CAM-NAV	2.73	1.52	1.47
6	C	800	HEM	CMD-C2D	2.69	1.55	1.47
6	A	800	HEM	C4D-ND	-2.68	1.34	1.39
2	A	2	RTZ	CAS-NAW	2.66	1.45	1.40
2	D	1	RTZ	CAM-NAV	2.66	1.52	1.47
2	C	2	RTZ	CAS-NAW	2.66	1.45	1.40
6	B	800	HEM	CHA-C4D	2.66	1.39	1.35
2	A	2	RTZ	CAL-CAN	2.65	1.57	1.52
6	C	800	HEM	FE-NB	2.62	2.07	1.97
2	A	2	RTZ	CAM-NAV	2.58	1.52	1.47
6	A	800	HEM	CHB-C1B	2.58	1.39	1.35
2	B	2	RTZ	CAR-NAW	2.57	1.45	1.40
2	C	2	RTZ	CAM-NAV	2.56	1.52	1.47
2	B	1	RTZ	CAR-NAW	2.54	1.45	1.40
2	D	1	RTZ	CAS-NAW	2.54	1.45	1.40
2	B	1	RTZ	CAM-NAV	2.54	1.52	1.47
2	B	1	RTZ	CAS-NAW	2.52	1.45	1.40
6	D	800	HEM	CMD-C2D	2.51	1.55	1.47
6	D	800	HEM	C4A-NA	-2.49	1.31	1.36
6	B	800	HEM	CHB-C1B	2.50	1.39	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	C	800	HEM	C4D-ND	-2.49	1.34	1.39
6	A	800	HEM	C4B-NB	-2.49	1.31	1.37
6	B	800	HEM	CMD-C2D	2.49	1.55	1.47
6	B	800	HEM	FE-NB	2.47	2.06	1.97
2	A	1	RTZ	CAS-NAW	2.46	1.44	1.40
6	D	800	HEM	FE-NB	2.46	2.06	1.97
2	B	2	RTZ	CAS-NAW	2.44	1.44	1.40
2	D	2	RTZ	CAS-NAW	2.43	1.44	1.40
2	D	1	RTZ	CAR-NAW	2.43	1.44	1.40
2	D	1	RTZ	CAH-CAR	2.41	1.44	1.39
6	A	800	HEM	CHA-C4D	2.40	1.39	1.35
6	C	800	HEM	O2A-CGA	-2.39	1.21	1.30
6	B	800	HEM	C4B-NB	-2.37	1.32	1.37
2	B	2	RTZ	CAO-CAQ	2.35	1.43	1.39
2	C	2	RTZ	CAO-CAQ	2.35	1.43	1.39
2	A	2	RTZ	CAO-CAQ	2.33	1.43	1.39
2	B	1	RTZ	CAO-CAS	2.33	1.43	1.39
2	C	1	RTZ	CAR-NAW	2.32	1.44	1.40
6	C	800	HEM	CBC-CAC	2.30	1.42	1.28
6	A	800	HEM	O2A-CGA	-2.29	1.22	1.30
2	D	2	RTZ	CAR-NAW	2.27	1.44	1.40
6	D	800	HEM	C4B-NB	-2.27	1.32	1.37
2	A	2	RTZ	CAO-CAS	2.25	1.43	1.39
6	D	800	HEM	CBC-CAC	2.25	1.41	1.28
2	C	1	RTZ	CAH-CAR	2.25	1.44	1.39
6	D	800	HEM	FE-NC	2.23	2.06	1.97
2	A	1	RTZ	CAH-CAR	2.22	1.43	1.39
6	B	800	HEM	C4A-NA	-2.22	1.32	1.36
2	A	1	RTZ	CAR-NAW	2.21	1.44	1.40
6	A	800	HEM	C4A-NA	-2.21	1.32	1.36
6	A	800	HEM	CBC-CAC	2.20	1.41	1.28
2	C	2	RTZ	CAO-CAS	2.20	1.43	1.39
6	C	800	HEM	CHA-C4D	2.19	1.39	1.35
6	B	800	HEM	O2A-CGA	-2.19	1.22	1.30
2	B	2	RTZ	CAH-CAR	2.19	1.43	1.39
6	B	800	HEM	CBC-CAC	2.17	1.41	1.28
2	D	2	RTZ	CAO-CAQ	2.17	1.43	1.39
2	D	2	RTZ	CAH-CAR	2.17	1.43	1.39
2	B	1	RTZ	CAO-CAQ	2.16	1.43	1.39
6	D	800	HEM	O2A-CGA	-2.15	1.22	1.30
2	B	1	RTZ	CAH-CAR	2.13	1.43	1.39
6	C	800	HEM	C4B-NB	-2.13	1.32	1.37

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	800	HEM	CHD-C4C	2.13	1.40	1.36
2	B	2	RTZ	CAO-CAS	2.12	1.43	1.39
6	C	800	HEM	C4A-NA	-2.10	1.32	1.36
2	C	2	RTZ	CAU-SAY	2.09	1.80	1.76
2	C	1	RTZ	CAS-NAW	2.08	1.44	1.40
2	C	2	RTZ	CAC-CAH	2.08	1.43	1.39
6	D	800	HEM	CHA-C4D	2.06	1.38	1.35
2	A	2	RTZ	CAC-CAH	2.03	1.43	1.39
2	D	2	RTZ	CAO-CAS	2.02	1.43	1.39
2	C	1	RTZ	CAU-SAY	2.01	1.80	1.76

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	2	RTZ	CAL-CAN-NAW	6.55	122.64	111.77
2	C	2	RTZ	CAL-CAN-NAW	6.44	122.45	111.77
2	A	1	RTZ	CAL-CAN-NAW	6.06	121.82	111.77
2	C	1	RTZ	CAL-CAN-NAW	5.93	121.61	111.77
2	D	1	RTZ	CAL-CAN-NAW	5.35	120.65	111.77
2	B	1	RTZ	CAL-CAN-NAW	5.29	120.55	111.77
2	A	2	RTZ	CAN-NAW-CAS	4.77	125.16	118.96
2	C	2	RTZ	CAN-NAW-CAS	4.73	125.12	118.96
6	B	800	HEM	CHD-C4C-NC	-4.44	120.87	124.73
6	A	800	HEM	CHD-C4C-NC	-4.40	120.91	124.73
2	C	2	RTZ	CAA-NAV-CAM	-4.33	102.94	110.56
2	A	2	RTZ	CAA-NAV-CAM	-4.18	103.20	110.56
2	A	1	RTZ	CAA-NAV-CAM	-4.03	103.47	110.56
2	C	1	RTZ	CAA-NAV-CAM	-4.02	103.48	110.56
2	B	1	RTZ	CAA-NAV-CAM	-3.99	103.55	110.56
6	D	800	HEM	CHD-C4C-NC	-3.97	121.28	124.73
2	D	1	RTZ	CAA-NAV-CAM	-3.93	103.65	110.56
6	B	800	HEM	C3B-C4B-NB	-3.66	111.38	114.00
6	D	800	HEM	C3B-C4B-NB	-3.64	111.39	114.00
6	C	800	HEM	CHD-C4C-NC	-3.62	121.59	124.73
2	B	2	RTZ	CAL-CAN-NAW	3.50	117.58	111.77
6	B	800	HEM	CHC-C1C-NC	-3.45	121.74	124.73
2	C	1	RTZ	CAN-CAL-CAP	-3.40	107.85	114.40
6	C	800	HEM	C3B-C4B-NB	-3.37	111.59	114.00
2	D	2	RTZ	CAL-CAN-NAW	3.33	117.30	111.77
2	B	1	RTZ	CAN-CAL-CAP	-3.21	108.22	114.40
6	C	800	HEM	C4A-CHB-C1B	3.18	131.66	127.47
6	A	800	HEM	C4A-CHB-C1B	3.16	131.63	127.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1	RTZ	CAN-CAL-CAP	-3.11	108.41	114.40
6	D	800	HEM	C4D-ND-C1D	3.04	108.27	105.16
6	B	800	HEM	C4D-ND-C1D	3.02	108.25	105.16
6	A	800	HEM	C4D-ND-C1D	3.02	108.25	105.16
6	A	800	HEM	C3B-C4B-NB	-3.00	111.85	114.00
6	B	800	HEM	C3A-C4A-NA	2.93	111.62	109.41
6	D	800	HEM	CHC-C1C-NC	-2.91	122.20	124.73
6	D	800	HEM	C1B-NB-C4B	2.90	108.12	105.16
6	B	800	HEM	C1B-NB-C4B	2.85	108.08	105.16
6	D	800	HEM	C4A-CHB-C1B	2.79	131.14	127.47
6	C	800	HEM	C3A-C4A-NA	2.79	111.52	109.41
6	A	800	HEM	C3A-C4A-NA	2.76	111.50	109.41
6	C	800	HEM	C1B-NB-C4B	2.74	107.96	105.16
2	A	1	RTZ	CAN-NAW-CAS	2.73	122.51	118.96
6	C	800	HEM	C4D-ND-C1D	2.71	107.93	105.16
2	D	2	RTZ	CAA-NAV-CAP	2.69	115.25	111.54
6	B	800	HEM	C2D-C1D-ND	-2.67	109.77	112.93
6	D	800	HEM	C2D-C1D-ND	-2.66	109.79	112.93
2	D	1	RTZ	CAL-CAP-CAG	-2.64	107.94	112.63
6	C	800	HEM	CHC-C1C-NC	-2.64	122.44	124.73
2	D	2	RTZ	CAM-NAV-CAP	2.64	115.19	112.18
2	B	2	RTZ	CAA-NAV-CAM	-2.63	105.94	110.56
6	A	800	HEM	C2D-C1D-ND	-2.62	109.83	112.93
2	B	1	RTZ	CAL-CAP-CAG	-2.62	107.98	112.63
6	D	800	HEM	C3A-C4A-NA	2.59	111.37	109.41
6	B	800	HEM	C4A-CHB-C1B	2.59	130.88	127.47
2	A	1	RTZ	CAL-CAP-CAG	-2.59	108.03	112.63
2	B	2	RTZ	CAA-NAV-CAP	2.57	115.09	111.54
2	B	2	RTZ	CAM-NAV-CAP	2.55	115.09	112.18
6	C	800	HEM	C2D-C1D-ND	-2.55	109.92	112.93
6	A	800	HEM	CHC-C1C-NC	-2.54	122.52	124.73
6	A	800	HEM	C1B-NB-C4B	2.54	107.76	105.16
2	D	1	RTZ	CAN-CAL-CAP	-2.53	109.52	114.40
2	D	2	RTZ	CAA-NAV-CAM	-2.52	106.13	110.56
2	C	1	RTZ	CAN-NAW-CAS	2.49	122.20	118.96
2	C	1	RTZ	CAL-CAP-CAG	-2.45	108.28	112.63
2	B	1	RTZ	CAN-NAW-CAS	2.34	122.00	118.96
2	D	1	RTZ	CAN-NAW-CAS	2.30	121.95	118.96
2	C	2	RTZ	CAS-NAW-CAR	-2.23	114.75	119.99
6	C	800	HEM	CHC-C4B-NB	-2.13	122.81	124.58
2	D	2	RTZ	CAN-NAW-CAS	2.12	121.72	118.96
6	C	800	HEM	CBA-CAA-C2A	2.09	116.38	112.69

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
6	A	800	HEM	CHC-C4B-NB	-2.07	122.86	124.58
6	B	800	HEM	CHC-C4B-NB	-2.04	122.89	124.58
6	B	800	HEM	C4C-NC-C1C	2.03	107.64	105.53

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

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

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	467/479 (97%)	0.10	17 (3%) 41 45	15, 29, 53, 80	0
1	B	456/479 (95%)	0.17	25 (5%) 24 26	17, 34, 62, 82	0
1	C	467/479 (97%)	0.15	21 (4%) 32 35	15, 30, 57, 94	0
1	D	456/479 (95%)	0.18	21 (4%) 31 34	18, 34, 57, 82	0
All	All	1846/1916 (96%)	0.15	84 (4%) 32 34	15, 32, 58, 94	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	43	LEU	7.6
1	A	38	LEU	7.3
1	C	41	PRO	7.2
1	C	48	HIS	7.0
1	C	47	LEU	6.9
1	A	40	LEU	6.8
1	A	43	LEU	6.3
1	D	39	PRO	6.1
1	C	40	LEU	6.1
1	C	46	LEU	6.0
1	B	75	TRP	6.0
1	C	38	LEU	6.0
1	B	38	LEU	5.9
1	A	47	LEU	5.8
1	D	65	PHE	5.8
1	B	39	PRO	5.7
1	A	41	PRO	5.4
1	A	48	HIS	5.3
1	D	38	LEU	5.1
1	A	75	TRP	4.8
1	B	145	GLY	4.5

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Mol	Chain	Res	Type	RSRZ
1	B	144	LEU	4.3
1	D	219	PHE	4.3
1	D	145	GLY	4.3
1	D	75	TRP	4.1
1	B	50	ASP	4.1
1	C	49	VAL	4.0
1	D	144	LEU	3.9
1	C	145	GLY	3.9
1	B	424	GLN	3.8
1	C	219	PHE	3.7
1	A	32	LYS	3.7
1	C	39	PRO	3.7
1	D	423	ALA	3.7
1	D	63	ARG	3.6
1	B	63	ARG	3.5
1	B	33	LEU	3.4
1	B	65	PHE	3.4
1	B	74	ALA	3.4
1	D	424	GLN	3.3
1	A	219	PHE	3.3
1	C	32	LYS	3.3
1	B	64	ARG	3.2
1	A	39	PRO	3.1
1	D	342	VAL	3.1
1	D	33	LEU	3.0
1	B	219	PHE	3.0
1	C	42	GLY	3.0
1	B	73	LEU	3.0
1	A	46	LEU	3.0
1	C	33	LEU	3.0
1	A	31	GLY	2.9
1	D	64	ARG	2.9
1	B	471	GLY	2.8
1	A	342	VAL	2.8
1	C	235	ALA	2.8
1	D	387	PHE	2.8
1	B	32	LYS	2.7
1	B	342	VAL	2.7
1	D	169	GLY	2.7
1	A	144	LEU	2.7
1	C	44	GLY	2.6
1	A	33	LEU	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	75	TRP	2.6
1	C	31	GLY	2.5
1	D	74	ALA	2.4
1	D	386	GLY	2.4
1	A	308	VAL	2.4
1	B	423	ALA	2.4
1	D	340	GLY	2.3
1	D	50	ASP	2.3
1	D	146	LYS	2.3
1	C	144	LEU	2.3
1	B	381	ASP	2.2
1	C	51	PHE	2.2
1	B	480	VAL	2.2
1	B	386	GLY	2.1
1	B	470	THR	2.1
1	A	484	LEU	2.1
1	B	387	PHE	2.1
1	D	168	SER	2.1
1	C	56	TYR	2.1
1	B	72	GLN	2.0
1	B	53	ASN	2.0

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

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

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(Å ²)	Q<0.9
4	GOL	A	752	6/6	0.32	24.90	55,58,59,60	0

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Mol	Type	Chain	Res	Atoms	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	GOL	B	750	6/6	0.24	6.89	42,51,52,55	0
4	GOL	D	750	6/6	0.18	4.47	46,47,48,49	0
4	GOL	A	751	6/6	0.22	2.94	70,71,71,72	0
4	GOL	A	753	6/6	0.22	2.88	31,42,46,47	0
4	GOL	C	750	6/6	0.13	1.95	33,36,37,38	0
4	GOL	A	750	6/6	0.13	1.71	34,41,42,45	0
5	PO4	A	790	5/5	0.24	1.70	109,109,110,110	0
2	RTZ	C	2	25/25	0.21	1.62	51,62,64,64	0
2	RTZ	D	2	25/25	0.15	1.17	48,50,52,55	0
2	RTZ	A	2	25/25	0.17	1.06	46,57,63,63	0
2	RTZ	B	2	25/25	0.17	0.82	53,54,55,57	0
2	RTZ	B	1	25/25	0.13	0.62	28,34,37,41	0
2	RTZ	D	1	25/25	0.14	0.53	29,35,38,39	0
2	RTZ	A	1	25/25	0.16	0.47	26,29,33,36	0
6	HEM	C	800	43/43	0.15	0.30	9,14,18,21	0
2	RTZ	C	1	25/25	0.14	0.29	25,29,32,35	0
6	HEM	D	800	43/43	0.14	0.23	16,22,23,24	0
6	HEM	A	800	43/43	0.15	0.21	9,14,17,19	0
6	HEM	B	800	43/43	0.13	-0.12	10,17,20,22	0
3	ZN	A	602	1/1	0.10	-0.88	90,90,90,90	0
3	ZN	C	602	1/1	0.08	-0.94	79,79,79,79	0
3	ZN	A	600	1/1	0.08	-1.05	21,21,21,21	0
3	ZN	D	600	1/1	0.07	-1.21	26,26,26,26	0
3	ZN	B	600	1/1	0.07	-1.45	23,23,23,23	0
3	ZN	C	600	1/1	0.07	-1.52	24,24,24,24	0
3	ZN	A	601	1/1	0.04	-1.55	27,27,27,27	0

6.5 Other polymers ⓘ

There are no such residues in this entry.