



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 06:31 AM GMT

PDB ID : 2XCI
Title : Membrane-embedded monofunctional glycosyltransferase WaaA of Aquifex aeolicus, substrate-free form
Authors : Schmidt, H.; Hansen, G.; Hilgenfeld, R.; Mamat, U.; Mesters, J.R.
Deposited on : 2010-04-26
Resolution : 2.00 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

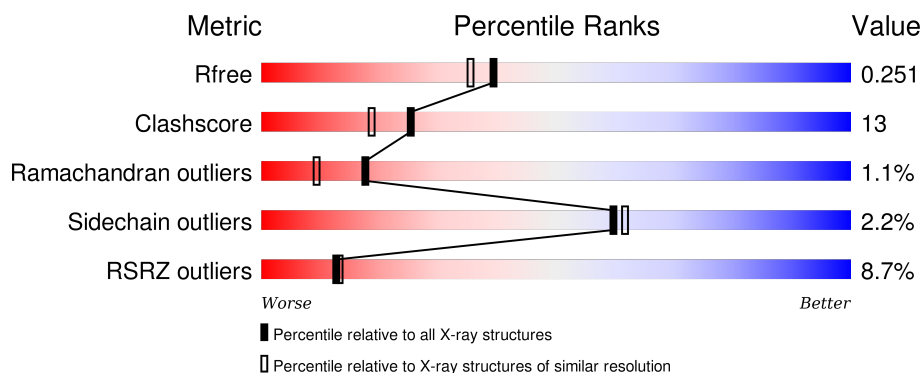
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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}	91344	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	374	<div> <div>7%</div> <div>70% 23% • 5%</div> </div>
1	B	374	<div> <div>6%</div> <div>73% 19% • 6%</div> </div>
1	C	374	<div> <div>11%</div> <div>70% 22% • 6%</div> </div>
1	D	374	<div> <div>8%</div> <div>78% 15% • 6%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	PG4	A	1353	-	-	-	X
2	PG4	A	1354	-	-	-	X
2	PG4	D	1353	-	-	X	X
2	PG4	D	1354	-	-	-	X
3	GOL	A	1356	-	-	X	X
3	GOL	A	1357	-	-	X	-
3	GOL	B	1353	-	-	-	X
3	GOL	C	1353	-	-	X	-
3	GOL	C	1354	-	-	X	X
3	GOL	D	1355	-	-	-	X
5	BME	C	1356	-	-	X	X
6	CL	D	1359	-	-	-	X

2 Entry composition

There are 7 unique types of molecules in this entry. The entry contains 11857 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 3-DEOXY-D-MANNO-2-OCTULOSONIC ACID TRANSFERASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	354	Total	C	N	O	S	0	0	0
			2888	1899	482	499	8			
1	B	352	Total	C	N	O	S	0	0	0
			2867	1887	475	497	8			
1	C	353	Total	C	N	O	S	0	0	0
			2877	1893	478	498	8			
1	D	352	Total	C	N	O	S	0	0	0
			2867	1887	475	497	8			

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-20	MET	-	EXPRESSION TAG	UNP O66663
A	-19	GLY	-	EXPRESSION TAG	UNP O66663
A	-18	HIS	-	EXPRESSION TAG	UNP O66663
A	-17	HIS	-	EXPRESSION TAG	UNP O66663
A	-16	HIS	-	EXPRESSION TAG	UNP O66663
A	-15	HIS	-	EXPRESSION TAG	UNP O66663
A	-14	HIS	-	EXPRESSION TAG	UNP O66663
A	-13	HIS	-	EXPRESSION TAG	UNP O66663
A	-12	HIS	-	EXPRESSION TAG	UNP O66663
A	-11	HIS	-	EXPRESSION TAG	UNP O66663
A	-10	HIS	-	EXPRESSION TAG	UNP O66663
A	-9	HIS	-	EXPRESSION TAG	UNP O66663
A	-8	SER	-	EXPRESSION TAG	UNP O66663
A	-7	SER	-	EXPRESSION TAG	UNP O66663
A	-6	GLY	-	EXPRESSION TAG	UNP O66663
A	-5	HIS	-	EXPRESSION TAG	UNP O66663
A	-4	ILE	-	EXPRESSION TAG	UNP O66663
A	-3	GLU	-	EXPRESSION TAG	UNP O66663
A	-2	GLY	-	EXPRESSION TAG	UNP O66663
A	-1	ARG	-	EXPRESSION TAG	UNP O66663

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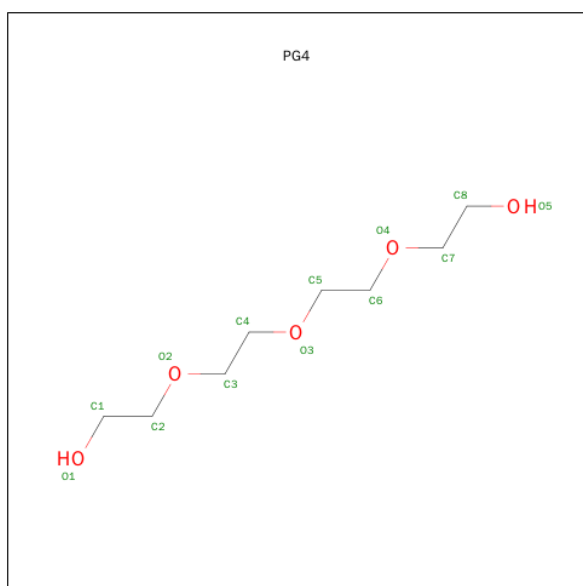
Chain	Residue	Modelled	Actual	Comment	Reference
A	0	HIS	-	EXPRESSION TAG	UNP O66663
B	-20	MET	-	EXPRESSION TAG	UNP O66663
B	-19	GLY	-	EXPRESSION TAG	UNP O66663
B	-18	HIS	-	EXPRESSION TAG	UNP O66663
B	-17	HIS	-	EXPRESSION TAG	UNP O66663
B	-16	HIS	-	EXPRESSION TAG	UNP O66663
B	-15	HIS	-	EXPRESSION TAG	UNP O66663
B	-14	HIS	-	EXPRESSION TAG	UNP O66663
B	-13	HIS	-	EXPRESSION TAG	UNP O66663
B	-12	HIS	-	EXPRESSION TAG	UNP O66663
B	-11	HIS	-	EXPRESSION TAG	UNP O66663
B	-10	HIS	-	EXPRESSION TAG	UNP O66663
B	-9	HIS	-	EXPRESSION TAG	UNP O66663
B	-8	SER	-	EXPRESSION TAG	UNP O66663
B	-7	SER	-	EXPRESSION TAG	UNP O66663
B	-6	GLY	-	EXPRESSION TAG	UNP O66663
B	-5	HIS	-	EXPRESSION TAG	UNP O66663
B	-4	ILE	-	EXPRESSION TAG	UNP O66663
B	-3	GLU	-	EXPRESSION TAG	UNP O66663
B	-2	GLY	-	EXPRESSION TAG	UNP O66663
B	-1	ARG	-	EXPRESSION TAG	UNP O66663
B	0	HIS	-	EXPRESSION TAG	UNP O66663
C	-20	MET	-	EXPRESSION TAG	UNP O66663
C	-19	GLY	-	EXPRESSION TAG	UNP O66663
C	-18	HIS	-	EXPRESSION TAG	UNP O66663
C	-17	HIS	-	EXPRESSION TAG	UNP O66663
C	-16	HIS	-	EXPRESSION TAG	UNP O66663
C	-15	HIS	-	EXPRESSION TAG	UNP O66663
C	-14	HIS	-	EXPRESSION TAG	UNP O66663
C	-13	HIS	-	EXPRESSION TAG	UNP O66663
C	-12	HIS	-	EXPRESSION TAG	UNP O66663
C	-11	HIS	-	EXPRESSION TAG	UNP O66663
C	-10	HIS	-	EXPRESSION TAG	UNP O66663
C	-9	HIS	-	EXPRESSION TAG	UNP O66663
C	-8	SER	-	EXPRESSION TAG	UNP O66663
C	-7	SER	-	EXPRESSION TAG	UNP O66663
C	-6	GLY	-	EXPRESSION TAG	UNP O66663
C	-5	HIS	-	EXPRESSION TAG	UNP O66663
C	-4	ILE	-	EXPRESSION TAG	UNP O66663
C	-3	GLU	-	EXPRESSION TAG	UNP O66663
C	-2	GLY	-	EXPRESSION TAG	UNP O66663
C	-1	ARG	-	EXPRESSION TAG	UNP O66663

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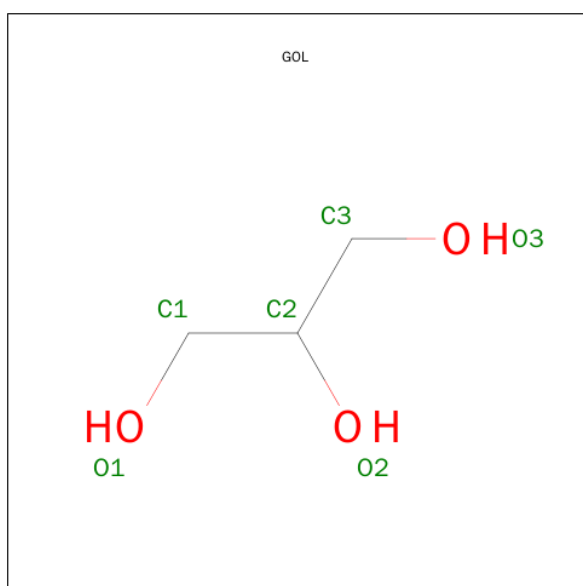
Chain	Residue	Modelled	Actual	Comment	Reference
C	0	HIS	-	EXPRESSION TAG	UNP O66663
D	-20	MET	-	EXPRESSION TAG	UNP O66663
D	-19	GLY	-	EXPRESSION TAG	UNP O66663
D	-18	HIS	-	EXPRESSION TAG	UNP O66663
D	-17	HIS	-	EXPRESSION TAG	UNP O66663
D	-16	HIS	-	EXPRESSION TAG	UNP O66663
D	-15	HIS	-	EXPRESSION TAG	UNP O66663
D	-14	HIS	-	EXPRESSION TAG	UNP O66663
D	-13	HIS	-	EXPRESSION TAG	UNP O66663
D	-12	HIS	-	EXPRESSION TAG	UNP O66663
D	-11	HIS	-	EXPRESSION TAG	UNP O66663
D	-10	HIS	-	EXPRESSION TAG	UNP O66663
D	-9	HIS	-	EXPRESSION TAG	UNP O66663
D	-8	SER	-	EXPRESSION TAG	UNP O66663
D	-7	SER	-	EXPRESSION TAG	UNP O66663
D	-6	GLY	-	EXPRESSION TAG	UNP O66663
D	-5	HIS	-	EXPRESSION TAG	UNP O66663
D	-4	ILE	-	EXPRESSION TAG	UNP O66663
D	-3	GLU	-	EXPRESSION TAG	UNP O66663
D	-2	GLY	-	EXPRESSION TAG	UNP O66663
D	-1	ARG	-	EXPRESSION TAG	UNP O66663
D	0	HIS	-	EXPRESSION TAG	UNP O66663

- Molecule 2 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			7	4	3		
2	A	1	Total	C	O	0	0
			12	8	4		
2	D	1	Total	C	O	0	0
			13	8	5		
2	D	1	Total	C	O	0	0
			10	6	4		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).

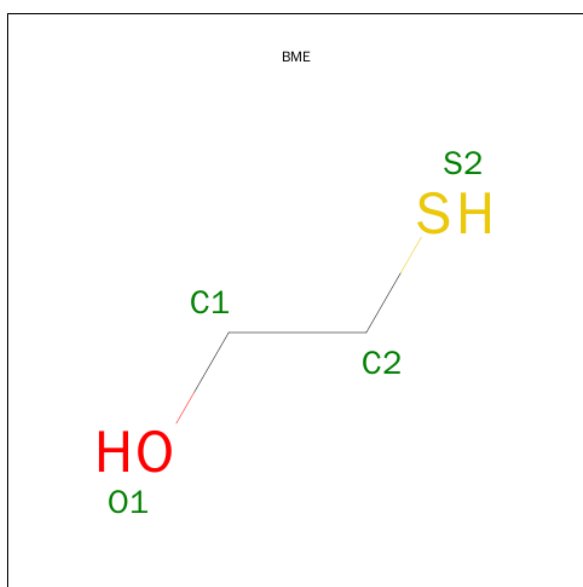


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	C	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		
3	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	B	1	Total Ni 1 1	0	0
4	A	1	Total Ni 1 1	0	0
4	D	1	Total Ni 1 1	0	0
4	C	1	Total Ni 1 1	0	0

- Molecule 5 is BETA-MERCAPTOETHANOL (three-letter code: BME) (formula: C₂H₆OS).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	C	1	Total C O S 4 2 1 1	0	0
5	C	1	Total C O S 4 2 1 1	0	0

- Molecule 6 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	B	2	Total Cl 2 2	0	0
6	A	3	Total Cl 3 3	0	0
6	D	2	Total Cl 2 2	0	0

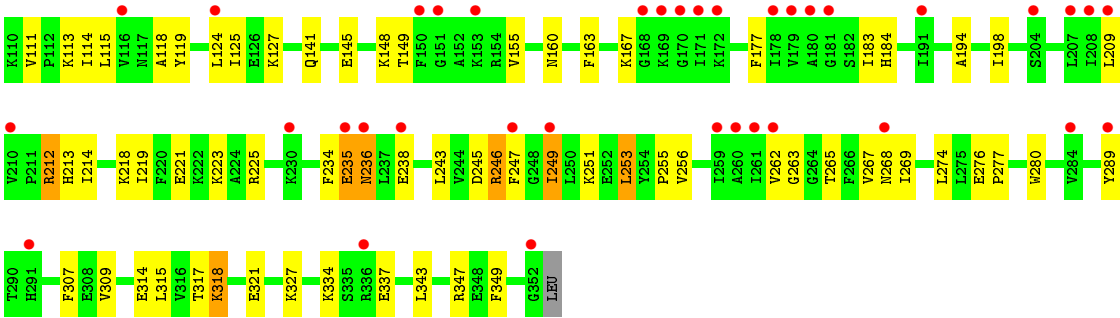
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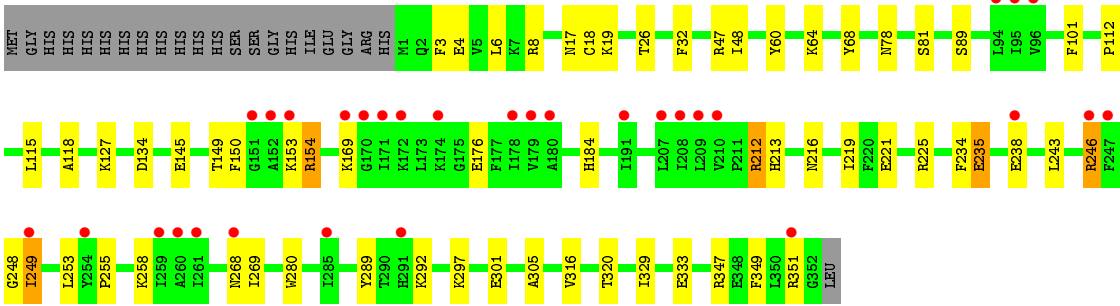
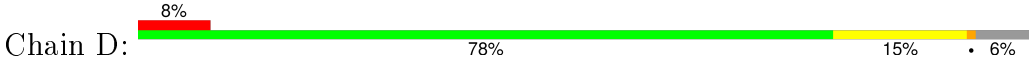
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	C	1	Total	Cl	0	0
			1	1		

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	57	Total	O	0	0
			57	57		
7	B	69	Total	O	0	0
			69	69		
7	C	58	Total	O	0	0
			58	58		
7	D	64	Total	O	0	0
			64	64		



● Molecule 1: 3-DEOXY-D-MANNO-2-OCTULOSONIC ACID TRANSFERASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	132.50Å 45.82Å 144.07Å 90.00° 97.23° 90.00°	Depositor
Resolution (Å)	29.85 – 2.00 29.85 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.9 (29.85-2.00) 98.9 (29.85-2.00)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 2.00Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.207 , 0.254 0.206 , 0.251	Depositor DCC
R_{free} test set	5821 reflections (5.28%)	DCC
Wilson B-factor (Å ²)	38.0	Xtriage
Anisotropy	0.712	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 55.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	0 of 115974 reflections	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	11857	wwPDB-VP
Average B, all atoms (Å ²)	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.23% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: GOL, PG4, CL, NI, BME

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	0.40	0/2952	0.55	0/3960
1	B	0.41	0/2930	0.55	0/3931
1	C	0.39	0/2941	0.55	0/3946
1	D	0.39	0/2930	0.54	0/3931
All	All	0.39	0/11753	0.55	0/15768

There are no bond length outliers.

There are no bond angle outliers.

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	2888	0	3028	87	0
1	B	2867	0	3008	75	0
1	C	2877	0	3015	84	0
1	D	2867	0	3008	52	0
2	A	19	0	24	3	0
2	D	23	0	31	10	0
3	A	18	0	24	10	0
3	B	6	0	8	3	0
3	C	12	0	16	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	D	12	0	16	4	0
4	A	1	0	0	0	0
4	B	1	0	0	0	0
4	C	1	0	0	0	0
4	D	1	0	0	0	0
5	C	8	0	12	8	0
6	A	3	0	0	0	0
6	B	2	0	0	0	0
6	C	1	0	0	0	0
6	D	2	0	0	0	0
7	A	57	0	0	2	0
7	B	69	0	0	1	0
7	C	58	0	0	2	0
7	D	64	0	0	5	0
All	All	11857	0	12190	299	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 13.

All (299) 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:221:GLU:HG2	1:B:231:THR:HG21	1.37	1.03
3:A:1357:GOL:H12	5:C:1356:BME:S2	2.02	0.99
3:B:1353:GOL:H11	7:B:2004:HOH:O	1.73	0.88
1:B:5:VAL:HG21	2:D:1353:PG4:H81	1.56	0.88
3:D:1355:GOL:H11	7:D:2001:HOH:O	1.75	0.85
1:B:176:GLU:HG2	1:B:205:LEU:HD23	1.56	0.85
1:D:219:ILE:H	1:D:219:ILE:HD12	1.41	0.85
1:C:6:LEU:HB2	3:C:1354:GOL:H32	1.56	0.85
1:A:185:THR:H	3:A:1357:GOL:H32	1.41	0.84
1:A:-1:ARG:HD2	1:C:119:TYR:OH	1.77	0.84
1:A:219:ILE:H	1:A:219:ILE:HD12	1.42	0.84
1:C:209:LEU:HD23	1:C:243:LEU:HD21	1.61	0.83
1:B:248:GLY:O	1:B:249:ILE:HG22	1.77	0.83
1:A:6:LEU:H	3:A:1356:GOL:H2	1.42	0.82
1:D:184:HIS:CE1	1:D:269:ILE:HD13	2.17	0.80
1:D:6:LEU:H	3:D:1355:GOL:H2	1.46	0.80
1:B:19:LYS:HE3	1:B:90:LYS:HD2	1.65	0.79
1:D:216:ASN:HA	1:D:219:ILE:HD13	1.63	0.79
1:B:5:VAL:HG22	2:D:1353:PG4:H42	1.64	0.78

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:153:LYS:O	1:D:154:ARG:HB2	1.82	0.78
1:A:216:ASN:HA	1:A:219:ILE:HD13	1.65	0.76
1:B:246:ARG:HH22	1:B:249:ILE:HD13	1.51	0.76
1:B:5:VAL:CG2	2:D:1353:PG4:H81	2.15	0.76
1:A:212:ARG:C	1:A:212:ARG:HD3	2.06	0.74
1:C:307:PHE:CD1	1:C:318:LYS:HE2	2.22	0.74
3:A:1356:GOL:H11	1:C:125:ILE:HB	1.70	0.73
1:A:297:LYS:HE2	1:A:301:GLU:OE2	1.89	0.72
1:B:6:LEU:H	3:B:1353:GOL:H2	1.56	0.71
1:A:262:VAL:HG21	1:A:274:LEU:HD22	1.72	0.70
1:C:6:LEU:CB	3:C:1354:GOL:H32	2.20	0.70
1:B:246:ARG:HH11	1:B:246:ARG:H	1.39	0.70
1:C:6:LEU:HB2	3:C:1354:GOL:C3	2.22	0.69
1:C:262:VAL:HG21	1:C:274:LEU:HD23	1.73	0.69
1:B:109:THR:OG1	1:B:113:LYS:HE3	1.93	0.69
3:A:1357:GOL:C1	5:C:1356:BME:S2	2.80	0.69
1:A:212:ARG:HD3	1:A:212:ARG:O	1.93	0.69
1:A:116:VAL:HG13	1:A:342:TYR:CE1	2.28	0.69
1:A:116:VAL:HG13	1:A:342:TYR:HE1	1.59	0.68
1:A:125:ILE:HB	3:C:1354:GOL:H31	1.74	0.68
1:C:246:ARG:HH11	1:C:246:ARG:HB2	1.58	0.68
1:D:234:PHE:O	1:D:235:GLU:HB2	1.91	0.68
1:A:262:VAL:HG21	1:A:274:LEU:CD2	2.24	0.67
1:C:124:LEU:HA	1:C:127:LYS:HE3	1.77	0.67
1:B:25:HIS:NE2	1:B:100:GLU:OE1	2.28	0.67
1:B:153:LYS:O	1:B:154:ARG:HB2	1.94	0.66
1:C:255:PRO:HG3	1:C:280:TRP:CD2	2.30	0.66
1:C:262:VAL:HG21	1:C:274:LEU:CD2	2.25	0.65
1:D:18:CYS:HA	1:D:47:ARG:NH2	2.12	0.65
1:C:267:VAL:HG23	1:C:269:ILE:HG12	1.78	0.64
1:C:255:PRO:HG3	1:C:280:TRP:CE2	2.33	0.64
1:A:219:ILE:O	1:A:222:LYS:HG3	1.97	0.64
1:A:255:PRO:HG3	1:A:280:TRP:CE2	2.33	0.63
1:B:41:GLU:HG3	1:B:343:LEU:HD11	1.79	0.63
1:D:101:PHE:H	2:D:1353:PG4:H72	1.64	0.63
1:C:98:GLU:O	1:C:99:ARG:HB2	1.98	0.62
1:A:220:PHE:HB2	1:A:243:LEU:HD21	1.80	0.62
1:C:78:ASN:ND2	1:C:81:SER:H	1.97	0.62
1:B:209:LEU:HB3	1:B:243:LEU:HD23	1.81	0.62
1:B:293:VAL:HG22	1:B:296:LEU:HB3	1.80	0.61
1:C:99:ARG:O	1:C:99:ARG:HG2	2.00	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:234:PHE:O	1:C:235:GLU:HB2	1.99	0.61
1:A:-1:ARG:HD3	1:A:3:PHE:HE2	1.66	0.61
1:B:234:PHE:O	1:B:235:GLU:HG2	2.01	0.61
1:A:255:PRO:HG3	1:A:280:TRP:CD2	2.37	0.60
1:A:303:GLU:HG3	7:A:2055:HOH:O	2.00	0.59
1:A:218:LYS:O	1:A:222:LYS:HG2	2.02	0.59
1:D:219:ILE:N	1:D:219:ILE:HD12	2.13	0.59
1:B:25:HIS:CD2	1:B:100:GLU:OE1	2.56	0.59
1:B:254:TYR:N	1:B:255:PRO:HD2	2.18	0.59
1:D:219:ILE:H	1:D:219:ILE:CD1	2.13	0.59
1:C:221:GLU:HB2	1:C:243:LEU:HD12	1.84	0.59
1:C:4:GLU:HA	3:C:1354:GOL:H12	1.85	0.59
1:D:127:LYS:HG2	1:D:150:PHE:CE1	2.38	0.58
1:A:219:ILE:HG23	1:C:223:LYS:NZ	2.18	0.58
1:C:19:LYS:HA	1:C:89:SER:O	2.03	0.58
1:D:316:VAL:O	1:D:320:THR:HG23	2.03	0.58
3:A:1356:GOL:H11	1:C:125:ILE:CB	2.33	0.58
1:B:316:VAL:O	1:B:320:THR:HG23	2.04	0.58
3:A:1356:GOL:H11	1:C:125:ILE:CG2	2.34	0.57
1:D:213:HIS:ND1	3:D:1356:GOL:H31	2.19	0.57
1:D:154:ARG:HG3	7:D:2024:HOH:O	2.05	0.57
1:B:209:LEU:HD23	1:B:243:LEU:HD21	1.87	0.57
1:A:248:GLY:O	1:A:249:ILE:HB	2.04	0.57
1:C:213:HIS:ND1	5:C:1356:BME:H21	2.20	0.57
1:A:82:VAL:CG1	1:A:109:THR:HG23	2.35	0.57
1:A:234:PHE:O	1:A:235:GLU:HB2	2.05	0.57
1:C:219:ILE:HD12	7:C:2038:HOH:O	2.05	0.56
1:A:281:GLY:HA2	1:A:329:ILE:HB	1.88	0.56
1:A:262:VAL:CG2	1:A:274:LEU:HD22	2.36	0.56
1:B:209:LEU:HB3	1:B:243:LEU:CD2	2.36	0.56
1:D:248:GLY:O	1:D:249:ILE:HB	2.06	0.56
1:C:109:THR:OG1	1:C:113:LYS:HE3	2.05	0.56
1:A:184:HIS:CE1	1:A:269:ILE:HD12	2.40	0.56
1:A:121:LYS:NZ	3:C:1353:GOL:C3	2.68	0.56
1:B:267:VAL:HG23	1:B:269:ILE:HG12	1.88	0.55
1:C:209:LEU:HB3	1:C:243:LEU:HD23	1.88	0.55
1:B:221:GLU:HB2	1:B:243:LEU:HD12	1.87	0.55
1:D:184:HIS:CE1	1:D:269:ILE:CD1	2.89	0.55
1:C:274:LEU:O	1:C:277:PRO:HD2	2.07	0.55
1:D:297:LYS:HE2	1:D:301:GLU:OE2	2.06	0.55
1:A:153:LYS:O	1:A:154:ARG:HB2	2.07	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:LYS:NZ	3:C:1353:GOL:H31	2.21	0.54
1:C:7:LYS:HG2	1:C:12:LYS:HG3	1.88	0.54
3:C:1353:GOL:H2	3:C:1354:GOL:O2	2.08	0.54
1:A:234:PHE:CD2	1:A:235:GLU:N	2.76	0.54
1:A:184:HIS:HB3	3:A:1357:GOL:H31	1.89	0.54
1:B:141:GLN:O	1:B:145:GLU:HG2	2.06	0.54
1:B:19:LYS:HE3	1:B:90:LYS:CD	2.36	0.54
1:D:212:ARG:HD3	1:D:212:ARG:C	2.28	0.54
1:B:222:LYS:HG2	1:B:225:ARG:NH2	2.23	0.54
1:B:72:HIS:ND1	1:B:73:PRO:HD2	2.23	0.54
1:B:176:GLU:HG3	1:B:258:LYS:HD2	1.89	0.54
1:B:3:PHE:CD1	2:D:1353:PG4:H62	2.43	0.53
1:B:3:PHE:HB2	2:D:1353:PG4:H82	1.89	0.53
1:A:268:ASN:OD1	1:A:291:HIS:HE1	1.91	0.53
1:A:267:VAL:HB	5:C:1355:BME:H12	1.91	0.53
1:D:212:ARG:O	1:D:212:ARG:HD3	2.09	0.53
1:D:19:LYS:HA	1:D:89:SER:O	2.08	0.52
1:A:221:GLU:HB2	1:A:243:LEU:HD22	1.90	0.52
1:C:78:ASN:HD21	1:C:81:SER:H	1.58	0.52
1:C:213:HIS:ND1	5:C:1356:BME:C2	2.73	0.52
1:B:153:LYS:O	1:B:154:ARG:CB	2.58	0.52
1:C:268:ASN:HA	1:C:289:TYR:CD1	2.45	0.52
1:A:114:ILE:HD11	1:A:349:PHE:CD1	2.45	0.52
1:C:218:LYS:HG3	7:C:2039:HOH:O	2.09	0.51
1:C:214:ILE:HG13	1:C:245:ASP:HA	1.91	0.51
1:D:176:GLU:HG3	1:D:258:LYS:CE	2.40	0.51
1:D:255:PRO:HG3	1:D:280:TRP:CD2	2.45	0.51
1:D:18:CYS:HA	1:D:47:ARG:HH21	1.74	0.51
1:D:213:HIS:CE1	3:D:1356:GOL:H31	2.45	0.51
1:B:184:HIS:NE2	1:B:212:ARG:HG3	2.25	0.51
1:A:4:GLU:O	1:A:8:ARG:HG3	2.11	0.51
1:A:16:LYS:O	1:A:19:LYS:HG3	2.11	0.51
1:A:-1:ARG:HD3	1:A:3:PHE:CE2	2.46	0.51
1:B:83:LYS:O	1:B:87:GLU:HG2	2.11	0.51
1:D:255:PRO:HG3	1:D:280:TRP:CE2	2.45	0.50
2:A:1354:PG4:H52	1:C:5:VAL:HG22	1.93	0.50
1:A:309:VAL:HG11	1:A:315:LEU:HB2	1.94	0.50
1:B:78:ASN:ND2	1:B:81:SER:H	2.09	0.50
1:C:309:VAL:HG13	1:C:314:GLU:HB2	1.93	0.50
1:B:25:HIS:HE2	1:B:100:GLU:CD	2.15	0.50
1:D:351:ARG:HB3	7:D:2064:HOH:O	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:114:ILE:HD11	1:C:349:PHE:CE1	2.47	0.50
1:A:109:THR:OG1	1:A:113:LYS:HE3	2.12	0.50
1:C:184:HIS:HA	5:C:1356:BME:H11	1.93	0.50
1:A:26:THR:HG21	1:A:32:PHE:HA	1.93	0.49
1:C:25:HIS:CD2	1:C:100:GLU:HG3	2.47	0.49
1:A:351:ARG:HD2	7:A:2057:HOH:O	2.11	0.49
1:A:-1:ARG:HD2	1:C:119:TYR:HH	1.78	0.49
1:C:82:VAL:CG1	1:C:109:THR:HG23	2.41	0.49
1:C:141:GLN:O	1:C:145:GLU:HG2	2.13	0.49
1:A:267:VAL:HG23	1:A:269:ILE:HG13	1.93	0.49
1:D:101:PHE:H	2:D:1353:PG4:C7	2.26	0.49
1:B:5:VAL:N	3:B:1353:GOL:H2	2.27	0.48
1:C:209:LEU:HB3	1:C:243:LEU:CD2	2.43	0.48
1:B:246:ARG:HB2	1:B:246:ARG:CZ	2.43	0.48
1:C:114:ILE:HD11	1:C:349:PHE:CD1	2.47	0.48
1:C:234:PHE:O	1:C:235:GLU:CB	2.61	0.48
1:D:246:ARG:NH1	1:D:246:ARG:HB3	2.29	0.48
1:C:317:THR:O	1:C:321:GLU:HG3	2.14	0.48
1:A:172:LYS:O	1:A:173:LEU:HD23	2.14	0.48
1:C:145:GLU:O	1:C:149:THR:HG23	2.13	0.48
1:A:316:VAL:O	1:A:320:THR:HG23	2.14	0.48
1:A:184:HIS:CE1	1:A:212:ARG:HG3	2.48	0.47
1:D:246:ARG:HB3	1:D:246:ARG:CZ	2.44	0.47
1:D:234:PHE:O	1:D:235:GLU:CB	2.61	0.47
1:A:262:VAL:CG2	1:A:274:LEU:CD2	2.92	0.47
1:B:27:ALA:HB2	1:B:100:GLU:CD	2.35	0.47
1:C:51:THR:HA	1:C:72:HIS:O	2.15	0.47
1:C:309:VAL:HG11	1:C:315:LEU:HB2	1.95	0.47
1:C:4:GLU:O	1:C:8:ARG:HG3	2.14	0.47
1:B:253:LEU:O	1:B:256:VAL:HB	2.15	0.47
1:C:262:VAL:CG2	1:C:274:LEU:HD23	2.44	0.47
1:C:253:LEU:O	1:C:256:VAL:HB	2.15	0.47
1:C:109:THR:HB	1:C:111:VAL:HG12	1.97	0.47
1:B:97:VAL:HG12	1:B:98:GLU:CG	2.45	0.47
1:A:199:LYS:HE3	1:A:205:LEU:O	2.14	0.47
1:A:145:GLU:O	1:A:149:THR:HG23	2.14	0.47
1:B:265:THR:HG22	1:B:271:GLY:HA2	1.97	0.47
1:A:219:ILE:HG23	1:C:223:LYS:HZ1	1.78	0.47
1:C:167:LYS:O	1:C:251:LYS:HE2	2.15	0.47
1:A:254:TYR:N	1:A:255:PRO:CD	2.78	0.47
1:C:249:ILE:HG22	1:C:253:LEU:HD22	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:121:LYS:HZ2	3:C:1353:GOL:H31	1.80	0.47
1:A:274:LEU:O	1:A:277:PRO:HD2	2.15	0.47
1:B:19:LYS:HA	1:B:89:SER:O	2.14	0.46
3:A:1356:GOL:H11	1:C:125:ILE:HG22	1.97	0.46
1:A:262:VAL:O	1:A:262:VAL:HG23	2.16	0.46
1:A:121:LYS:NZ	3:C:1353:GOL:H32	2.31	0.46
1:C:20:GLY:O	1:C:92:LYS:HE2	2.16	0.46
1:B:293:VAL:CG2	1:B:296:LEU:HD23	2.46	0.46
1:B:267:VAL:HG23	1:B:269:ILE:CG1	2.46	0.46
1:C:343:LEU:HG	1:C:347:ARG:CZ	2.46	0.46
1:A:214:ILE:HG21	1:A:245:ASP:O	2.16	0.46
1:D:4:GLU:O	1:D:8:ARG:HG3	2.16	0.46
1:A:187:GLU:HG2	1:A:266:PHE:HB2	1.98	0.46
1:B:244:VAL:HG11	1:B:249:ILE:HG21	1.98	0.46
1:C:183:ILE:O	5:C:1356:BME:H11	2.16	0.46
3:A:1355:GOL:H11	3:A:1356:GOL:H12	1.98	0.46
1:B:187:GLU:HG2	1:B:267:VAL:HG13	1.98	0.46
1:D:268:ASN:HA	1:D:289:TYR:CG	2.51	0.46
1:D:221:GLU:HA	1:D:243:LEU:HD11	1.97	0.46
1:B:265:THR:CG2	1:B:271:GLY:HA2	2.46	0.45
1:B:221:GLU:N	1:B:243:LEU:HD11	2.32	0.45
1:B:312:GLU:O	1:B:316:VAL:HG23	2.17	0.45
1:B:198:ILE:HD12	1:B:316:VAL:HG13	1.98	0.45
1:A:234:PHE:O	1:A:235:GLU:CB	2.63	0.45
1:C:262:VAL:CG2	1:C:274:LEU:CD2	2.95	0.45
1:B:234:PHE:O	1:B:235:GLU:CB	2.63	0.45
1:D:249:ILE:HG23	1:D:253:LEU:CD2	2.46	0.45
1:D:60:TYR:CZ	1:D:64:LYS:HE3	2.52	0.45
1:C:212:ARG:HB3	5:C:1356:BME:H21	1.99	0.45
1:A:176:GLU:HG3	1:A:205:LEU:HA	1.98	0.45
1:C:115:LEU:HG	1:C:118:ALA:HB2	1.98	0.45
1:A:177:PHE:CD2	1:A:256:VAL:HG12	2.52	0.45
1:B:3:PHE:CB	2:D:1353:PG4:H82	2.47	0.45
1:A:219:ILE:H	1:A:219:ILE:CD1	2.16	0.45
1:C:3:PHE:HB2	3:C:1353:GOL:O3	2.17	0.45
1:C:263:GLY:O	1:C:265:THR:HG23	2.17	0.45
1:C:148:LYS:HG3	1:C:155:VAL:HG11	1.99	0.44
1:B:210:VAL:HG22	1:B:244:VAL:HG21	1.99	0.44
1:B:97:VAL:HG12	1:B:98:GLU:HG2	1.99	0.44
1:A:334:LYS:O	1:A:337:GLU:HB3	2.17	0.44
1:D:112:PRO:HA	1:D:134:ASP:OD2	2.16	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:153:LYS:O	1:D:154:ARG:CB	2.55	0.44
1:B:109:THR:OG1	1:B:113:LYS:CE	2.65	0.44
1:A:51:THR:HA	1:A:72:HIS:O	2.18	0.44
1:B:244:VAL:CG1	1:B:249:ILE:HG21	2.47	0.44
1:A:343:LEU:O	1:A:347:ARG:HG3	2.18	0.44
1:B:231:THR:HG22	1:B:232:SER:N	2.33	0.44
1:C:243:LEU:HD23	1:C:243:LEU:HA	1.65	0.44
1:B:246:ARG:HH22	1:B:249:ILE:CD1	2.25	0.44
1:B:177:PHE:CD1	1:B:206:LYS:HB2	2.53	0.44
2:A:1354:PG4:H52	1:C:5:VAL:CG2	2.48	0.43
1:A:217:ALA:O	1:A:243:LEU:CD2	2.66	0.43
1:B:53:PHE:CZ	2:D:1353:PG4:H21	2.53	0.43
1:D:176:GLU:HG3	1:D:258:LYS:HE2	2.01	0.43
1:A:294:ASN:O	1:A:298:GLU:HG3	2.18	0.43
1:B:51:THR:HA	1:B:72:HIS:O	2.19	0.43
1:A:16:LYS:C	1:A:18:CYS:H	2.22	0.43
1:B:274:LEU:O	1:B:277:PRO:HD2	2.18	0.43
1:A:167:LYS:O	1:A:251:LYS:HE2	2.19	0.43
1:A:78:ASN:ND2	1:A:81:SER:H	2.16	0.43
1:A:121:LYS:HZ3	3:C:1353:GOL:C3	2.31	0.43
1:C:148:LYS:HG3	1:C:155:VAL:CG1	2.48	0.43
1:A:184:HIS:CE1	1:A:269:ILE:CD1	3.01	0.43
1:D:17:ASN:O	1:D:47:ARG:NH2	2.52	0.42
1:D:221:GLU:HB2	1:D:243:LEU:HD13	2.01	0.42
1:D:48:ILE:HB	1:D:68:TYR:HA	2.01	0.42
1:C:184:HIS:NE2	1:C:212:ARG:HG3	2.33	0.42
1:C:194:ALA:O	1:C:198:ILE:HG12	2.19	0.42
1:A:121:LYS:HZ3	3:C:1353:GOL:H32	1.84	0.42
1:A:220:PHE:CB	1:A:243:LEU:HD21	2.47	0.42
1:D:47:ARG:HG3	1:D:47:ARG:HH11	1.83	0.42
1:B:258:LYS:HB3	1:B:258:LYS:HE2	1.77	0.42
1:C:262:VAL:HG21	1:C:274:LEU:HD22	2.00	0.42
1:A:153:LYS:O	1:A:154:ARG:CB	2.66	0.42
1:C:334:LYS:O	1:C:337:GLU:HB3	2.20	0.42
1:A:262:VAL:HG21	1:A:274:LEU:HD23	1.99	0.42
1:A:249:ILE:HG22	1:A:253:LEU:HD22	2.01	0.42
1:A:99:ARG:O	1:A:100:GLU:C	2.58	0.42
1:B:4:GLU:O	1:B:8:ARG:HG3	2.20	0.42
1:A:121:LYS:HZ2	3:C:1353:GOL:C3	2.32	0.42
1:B:234:PHE:O	1:B:235:GLU:CG	2.68	0.42
1:C:219:ILE:H	1:C:219:ILE:HD12	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:309:VAL:HG13	1:A:314:GLU:HB2	2.01	0.42
1:D:292:LYS:HD3	7:D:2046:HOH:O	2.19	0.42
1:A:218:LYS:HG2	1:A:222:LYS:NZ	2.35	0.41
1:B:255:PRO:HA	1:B:282:ILE:HD11	2.01	0.41
1:D:347:ARG:HG2	7:D:2006:HOH:O	2.19	0.41
1:D:78:ASN:ND2	1:D:81:SER:H	2.18	0.41
1:C:177:PHE:CD2	1:C:256:VAL:HG13	2.55	0.41
1:A:199:LYS:NZ	1:A:240:ASP:OD2	2.52	0.41
1:D:3:PHE:HD1	2:D:1354:PG4:H32	1.84	0.41
1:C:89:SER:O	1:C:90:LYS:C	2.58	0.41
1:C:7:LYS:HD3	1:C:12:LYS:CD	2.50	0.41
1:A:-1:ARG:C	1:A:1:MET:H	2.22	0.41
1:B:255:PRO:HA	1:B:282:ILE:CD1	2.50	0.41
1:B:221:GLU:CA	1:B:243:LEU:HD11	2.50	0.41
1:B:99:ARG:NH1	1:B:126:GLU:OE1	2.48	0.41
1:B:237:LEU:HA	1:B:242:ILE:HG12	2.03	0.41
1:B:237:LEU:HD23	1:B:237:LEU:N	2.36	0.41
1:C:236:ASN:HD22	1:C:236:ASN:HA	1.68	0.41
1:D:145:GLU:O	1:D:149:THR:HG23	2.21	0.41
1:A:219:ILE:HG23	1:C:223:LYS:HZ2	1.86	0.41
1:D:154:ARG:HH11	1:D:349:PHE:HD2	1.67	0.41
1:A:153:LYS:HE2	1:A:153:LYS:HB3	1.94	0.41
1:B:173:LEU:HD12	1:B:256:VAL:HG11	2.02	0.41
1:D:221:GLU:OE2	1:D:225:ARG:NH1	2.54	0.41
1:D:305:ALA:HB2	1:D:329:ILE:HD12	2.03	0.41
1:C:4:GLU:HA	3:C:1354:GOL:C1	2.51	0.41
1:C:160:ASN:HB3	1:C:163:PHE:CD2	2.56	0.40
1:C:236:ASN:HB3	1:C:238:GLU:OE1	2.21	0.40
1:D:115:LEU:HG	1:D:118:ALA:HB2	2.03	0.40
1:D:26:THR:HG21	1:D:32:PHE:HA	2.03	0.40
2:A:1354:PG4:H32	1:C:8:ARG:CZ	2.51	0.40
1:B:145:GLU:O	1:B:149:THR:HG23	2.21	0.40
1:A:114:ILE:HD11	1:A:349:PHE:CE1	2.56	0.40
1:B:270:GLY:HA2	1:B:291:HIS:CE1	2.55	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	352/374 (94%)	336 (96%)	11 (3%)	5 (1%)	14	6
1	B	350/374 (94%)	333 (95%)	12 (3%)	5 (1%)	14	6
1	C	351/374 (94%)	340 (97%)	8 (2%)	3 (1%)	21	13
1	D	350/374 (94%)	341 (97%)	6 (2%)	3 (1%)	21	13
All	All	1403/1496 (94%)	1350 (96%)	37 (3%)	16 (1%)	17	9

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	100	GLU
1	B	235	GLU
1	C	99	ARG
1	D	154	ARG
1	D	235	GLU
1	A	154	ARG
1	A	235	GLU
1	B	154	ARG
1	C	235	GLU
1	D	249	ILE
1	A	17	ASN
1	B	98	GLU
1	A	98	GLU
1	A	99	ARG
1	C	249	ILE
1	B	249	ILE

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	317/334 (95%)	312 (98%)	5 (2%)	70	73
1	B	315/334 (94%)	307 (98%)	8 (2%)	55	55
1	C	316/334 (95%)	306 (97%)	10 (3%)	46	44
1	D	315/334 (94%)	310 (98%)	5 (2%)	70	73
All	All	1263/1336 (94%)	1235 (98%)	28 (2%)	60	62

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	98	GLU
1	A	212	ARG
1	A	223	LYS
1	A	246	ARG
1	A	247	PHE
1	B	18	CYS
1	B	212	ARG
1	B	214	ILE
1	B	237	LEU
1	B	238	GLU
1	B	246	ARG
1	B	253	LEU
1	B	283	PRO
1	C	18	CYS
1	C	212	ARG
1	C	225	ARG
1	C	236	ASN
1	C	246	ARG
1	C	247	PHE
1	C	253	LEU
1	C	276	GLU
1	C	318	LYS
1	C	327	LYS
1	D	169	LYS
1	D	212	ARG
1	D	238	GLU
1	D	246	ARG
1	D	333	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (15) such

sidechains are listed below:

Mol	Chain	Res	Type
1	A	78	ASN
1	A	141	GLN
1	A	216	ASN
1	A	291	HIS
1	B	78	ASN
1	B	216	ASN
1	B	291	HIS
1	C	78	ASN
1	C	141	GLN
1	C	216	ASN
1	C	236	ASN
1	C	291	HIS
1	D	78	ASN
1	D	216	ASN
1	D	291	HIS

5.3.3 RNA ⓘ

There are no RNA molecules 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 26 ligands modelled in this entry, 12 are monoatomic - leaving 14 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	PG4	A	1353	-	6,6,12	0.51	0	5,5,11	1.31	0
2	PG4	A	1354	-	11,11,12	0.54	0	10,10,11	1.26	0
3	GOL	A	1355	-	5,5,5	0.34	0	5,5,5	0.54	0
3	GOL	A	1356	-	5,5,5	0.38	0	5,5,5	0.66	0
3	GOL	A	1357	-	5,5,5	0.37	0	5,5,5	0.43	0
3	GOL	B	1353	-	5,5,5	0.38	0	5,5,5	0.26	0
3	GOL	C	1353	-	5,5,5	0.34	0	5,5,5	0.41	0
3	GOL	C	1354	-	5,5,5	0.31	0	5,5,5	0.89	0
5	BME	C	1355	-	3,3,3	0.37	0	2,2,2	0.31	0
5	BME	C	1356	-	3,3,3	0.35	0	2,2,2	0.62	0
2	PG4	D	1353	-	12,12,12	0.50	0	11,11,11	1.80	6 (54%)
2	PG4	D	1354	-	9,9,12	0.51	0	8,8,11	1.42	0
3	GOL	D	1355	-	5,5,5	0.41	0	5,5,5	0.51	0
3	GOL	D	1356	-	5,5,5	0.55	0	5,5,5	0.43	0

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	PG4	A	1353	-	-	0/4/4/10	0/0/0/0
2	PG4	A	1354	-	-	0/9/9/10	0/0/0/0
3	GOL	A	1355	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1356	-	-	0/4/4/4	0/0/0/0
3	GOL	A	1357	-	-	0/4/4/4	0/0/0/0
3	GOL	B	1353	-	-	0/4/4/4	0/0/0/0
3	GOL	C	1353	-	-	0/4/4/4	0/0/0/0
3	GOL	C	1354	-	-	0/4/4/4	0/0/0/0
5	BME	C	1355	-	-	0/1/1/1	0/0/0/0
5	BME	C	1356	-	-	0/1/1/1	0/0/0/0
2	PG4	D	1353	-	-	0/10/10/10	0/0/0/0
2	PG4	D	1354	-	-	0/7/7/10	0/0/0/0
3	GOL	D	1355	-	-	0/4/4/4	0/0/0/0
3	GOL	D	1356	-	-	0/4/4/4	0/0/0/0

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1353	PG4	O3-C4-C3	2.01	119.31	110.36
2	D	1353	PG4	O4-C6-C5	2.02	119.32	110.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	1353	PG4	O4-C7-C8	2.08	120.02	110.43
2	D	1353	PG4	O3-C5-C6	2.15	119.92	110.36
2	D	1353	PG4	C5-O3-C4	2.21	122.81	113.31
2	D	1353	PG4	O2-C3-C4	2.22	120.23	110.36

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

13 monomers are involved in 51 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1354	PG4	3	0
3	A	1355	GOL	1	0
3	A	1356	GOL	6	0
3	A	1357	GOL	4	0
3	B	1353	GOL	3	0
3	C	1353	GOL	9	0
3	C	1354	GOL	7	0
5	C	1355	BME	1	0
5	C	1356	BME	7	0
2	D	1353	PG4	9	0
2	D	1354	PG4	1	0
3	D	1355	GOL	2	0
3	D	1356	GOL	2	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	354/374 (94%)	0.36	28 (7%) 15 16	28, 54, 85, 133	0
1	B	352/374 (94%)	0.39	24 (6%) 20 22	27, 56, 91, 117	0
1	C	353/374 (94%)	0.42	40 (11%) 7 7	28, 56, 90, 112	0
1	D	352/374 (94%)	0.40	31 (8%) 12 13	28, 56, 94, 123	0
All	All	1411/1496 (94%)	0.39	123 (8%) 13 13	27, 55, 91, 133	0

All (123) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	247	PHE	6.7
1	A	247	PHE	6.3
1	D	247	PHE	6.0
1	A	352	GLY	5.2
1	D	261	ILE	5.1
1	D	180	ALA	5.0
1	C	179	VAL	5.0
1	C	169	LYS	4.7
1	A	-1	ARG	4.6
1	C	247	PHE	4.6
1	D	179	VAL	4.5
1	B	169	LYS	4.5
1	A	261	ILE	4.4
1	C	178	ILE	4.4
1	C	261	ILE	4.4
1	A	179	VAL	4.3
1	C	289	TYR	4.2
1	B	153	LYS	4.1
1	D	208	ILE	4.1
1	D	178	ILE	4.0
1	C	180	ALA	3.8

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Mol	Chain	Res	Type	RSRZ
1	C	207	LEU	3.7
1	D	260	ALA	3.7
1	D	153	LYS	3.7
1	C	170	GLY	3.6
1	A	169	LYS	3.6
1	D	246	ARG	3.6
1	B	178	ILE	3.6
1	B	261	ILE	3.6
1	A	207	LEU	3.6
1	B	260	ALA	3.6
1	C	171	ILE	3.5
1	A	178	ILE	3.4
1	C	260	ALA	3.4
1	C	259	ILE	3.4
1	D	96	VAL	3.4
1	C	352	GLY	3.4
1	B	179	VAL	3.3
1	C	249	ILE	3.3
1	B	246	ARG	3.3
1	A	260	ALA	3.2
1	A	180	ALA	3.2
1	D	259	ILE	3.2
1	B	209	LEU	3.2
1	D	351	ARG	3.1
1	D	95	ILE	3.1
1	C	97	VAL	3.0
1	C	262	VAL	3.0
1	B	208	ILE	3.0
1	D	249	ILE	3.0
1	A	172	LYS	2.9
1	D	238	GLU	2.9
1	C	208	ILE	2.9
1	A	236	ASN	2.9
1	A	151	GLY	2.8
1	A	191	ILE	2.8
1	A	208	ILE	2.8
1	C	181	GLY	2.8
1	A	262	VAL	2.8
1	D	94	LEU	2.7
1	C	209	LEU	2.7
1	D	191	ILE	2.7
1	D	169	LYS	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	180	ALA	2.7
1	D	254	TYR	2.6
1	C	191	ILE	2.6
1	C	235	GLU	2.6
1	A	97	VAL	2.6
1	B	171	ILE	2.6
1	C	151	GLY	2.5
1	D	268	ASN	2.5
1	B	207	LEU	2.5
1	C	291	HIS	2.5
1	C	210	VAL	2.5
1	B	35	PHE	2.5
1	D	207	LEU	2.5
1	C	336	ARG	2.5
1	C	116	VAL	2.5
1	D	171	ILE	2.5
1	A	181	GLY	2.5
1	D	170	GLY	2.5
1	B	259	ILE	2.4
1	D	210	VAL	2.4
1	D	174	LYS	2.4
1	D	291	HIS	2.4
1	A	209	LEU	2.4
1	C	150	PHE	2.4
1	B	235	GLU	2.4
1	A	268	ASN	2.4
1	A	116	VAL	2.3
1	B	309	VAL	2.3
1	C	124	LEU	2.3
1	A	210	VAL	2.3
1	D	285	ILE	2.3
1	A	289	TYR	2.3
1	C	204	SER	2.3
1	D	151	GLY	2.3
1	B	96	VAL	2.3
1	A	259	ILE	2.3
1	A	285	ILE	2.3
1	C	24	VAL	2.2
1	A	154	ARG	2.2
1	B	116	VAL	2.2
1	A	148	LYS	2.2
1	B	268	ASN	2.2

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Mol	Chain	Res	Type	RSRZ
1	B	168	GLY	2.2
1	A	168	GLY	2.2
1	C	59	GLU	2.2
1	D	172	LYS	2.2
1	C	25	HIS	2.1
1	C	168	GLY	2.1
1	C	153	LYS	2.1
1	C	268	ASN	2.1
1	B	56	ARG	2.1
1	C	284	VAL	2.1
1	C	236	ASN	2.1
1	B	352	GLY	2.1
1	D	152	ALA	2.1
1	C	238	GLU	2.0
1	C	230	LYS	2.0
1	D	209	LEU	2.0
1	C	172	LYS	2.0
1	B	172	LYS	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 carbohydrates in this entry.

6.4 Ligands [i](#)

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	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
3	GOL	C	1354	6/6	0.70	0.27	6.63	55,61,65,66	0
2	PG4	A	1353	7/13	0.87	0.30	4.81	49,50,55,61	0
3	GOL	D	1355	6/6	0.87	0.20	4.28	49,55,59,59	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
3	GOL	B	1353	6/6	0.90	0.19	3.88	45,53,56,59	0
2	PG4	A	1354	12/13	0.85	0.25	3.83	41,58,69,71	0
5	BME	C	1356	4/4	0.84	0.23	3.52	44,45,51,67	0
6	CL	D	1359	1/1	0.69	0.17	3.20	74,74,74,74	0
3	GOL	A	1356	6/6	0.65	0.28	3.07	48,56,60,64	0
2	PG4	D	1354	10/13	0.82	0.23	2.87	51,57,71,72	0
2	PG4	D	1353	13/13	0.84	0.18	2.05	43,58,61,66	0
3	GOL	A	1355	6/6	0.85	0.18	1.68	42,51,66,71	0
3	GOL	A	1357	6/6	0.82	0.12	0.34	49,50,56,56	0
3	GOL	D	1356	6/6	0.79	0.13	-0.24	40,53,56,57	0
3	GOL	C	1353	6/6	0.84	0.11	-0.29	47,54,56,57	0
5	BME	C	1355	4/4	0.89	0.11	-0.70	66,67,67,68	0
6	CL	B	1355	1/1	0.96	0.08	-4.84	50,50,50,50	0
4	NI	B	1354	1/1	0.93	0.05	-	61,61,61,61	1
4	NI	D	1358	1/1	0.92	0.04	-	59,59,59,59	1
4	NI	C	1357	1/1	0.94	0.05	-	60,60,60,60	1
6	CL	A	1361	1/1	0.92	0.27	-	77,77,77,77	0
6	CL	A	1359	1/1	0.82	0.33	-	84,84,84,84	0
6	CL	D	1357	1/1	0.91	0.23	-	80,80,80,80	0
6	CL	B	1356	1/1	0.76	0.30	-	93,93,93,93	0
6	CL	C	1358	1/1	0.88	0.31	-	93,93,93,93	0
4	NI	A	1358	1/1	0.95	0.04	-	64,64,64,64	1
6	CL	A	1360	1/1	0.98	0.26	-	86,86,86,86	0

6.5 Other polymers [i](#)

There are no such residues in this entry.