



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:14 AM GMT

PDB ID : 3LCW
Title : L-KDO aldolase complexed with hydroxypyruvate
Authors : Chou, C.-Y.; Wang, A.H.-J.; Ko, T.-P.
Deposited on : 2010-01-12
Resolution : 2.35 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

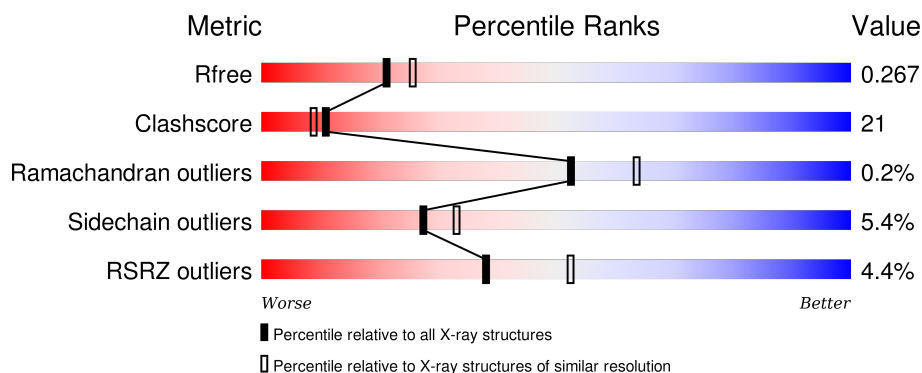
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.35 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	1352 (2.38-2.34)
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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	319	<div> <div>2%</div> <div>59%</div> <div>30%</div> <div>9%</div> </div>
1	B	319	<div> <div>6%</div> <div>62%</div> <div>26%</div> <div>9%</div> </div>
1	C	319	<div> <div>7%</div> <div>57%</div> <div>32%</div> <div>9%</div> </div>
1	D	319	<div> <div>2%</div> <div>56%</div> <div>32%</div> <div>9%</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	3PY	A	640	-	-	-	X
2	3PY	B	641	-	-	-	X
2	3PY	C	642	-	-	-	X
2	3PY	D	643	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 9473 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 N-acetylneuraminase lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	291	Total	C	N	O	S	71	0	0
			2234	1425	380	418	11			
1	B	291	Total	C	N	O	S	70	0	0
			2234	1425	380	418	11			
1	C	291	Total	C	N	O	S	68	0	0
			2234	1425	380	418	11			
1	D	291	Total	C	N	O	S	54	0	0
			2234	1425	380	418	11			

There are 120 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-21	MET	-	EXPRESSION TAG	UNP P0A6L4
A	-20	GLY	-	EXPRESSION TAG	UNP P0A6L4
A	-19	HIS	-	EXPRESSION TAG	UNP P0A6L4
A	-18	HIS	-	EXPRESSION TAG	UNP P0A6L4
A	-17	HIS	-	EXPRESSION TAG	UNP P0A6L4
A	-16	HIS	-	EXPRESSION TAG	UNP P0A6L4
A	-15	HIS	-	EXPRESSION TAG	UNP P0A6L4
A	-14	HIS	-	EXPRESSION TAG	UNP P0A6L4
A	-13	HIS	-	EXPRESSION TAG	UNP P0A6L4
A	-12	HIS	-	EXPRESSION TAG	UNP P0A6L4
A	-11	HIS	-	EXPRESSION TAG	UNP P0A6L4
A	-10	HIS	-	EXPRESSION TAG	UNP P0A6L4
A	-9	SER	-	EXPRESSION TAG	UNP P0A6L4
A	-8	SER	-	EXPRESSION TAG	UNP P0A6L4
A	-7	GLY	-	EXPRESSION TAG	UNP P0A6L4
A	-6	HIS	-	EXPRESSION TAG	UNP P0A6L4
A	-5	ILE	-	EXPRESSION TAG	UNP P0A6L4
A	-4	GLU	-	EXPRESSION TAG	UNP P0A6L4
A	-3	GLY	-	EXPRESSION TAG	UNP P0A6L4
A	-2	ARG	-	EXPRESSION TAG	UNP P0A6L4
A	-1	HIS	-	EXPRESSION TAG	UNP P0A6L4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MET	-	EXPRESSION TAG	UNP P0A6L4
A	60	ALA	GLU	ENGINEERED MUTATION	UNP P0A6L4
A	98	HIS	TYR	ENGINEERED MUTATION	UNP P0A6L4
A	115	LEU	PHE	ENGINEERED MUTATION	UNP P0A6L4
A	150	GLY	ASP	ENGINEERED MUTATION	UNP P0A6L4
A	153	TYR	ASN	ENGINEERED MUTATION	UNP P0A6L4
A	251	ILE	VAL	ENGINEERED MUTATION	UNP P0A6L4
A	265	ILE	VAL	ENGINEERED MUTATION	UNP P0A6L4
A	281	CYS	TYR	ENGINEERED MUTATION	UNP P0A6L4
B	-21	MET	-	EXPRESSION TAG	UNP P0A6L4
B	-20	GLY	-	EXPRESSION TAG	UNP P0A6L4
B	-19	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	-18	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	-17	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	-16	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	-15	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	-14	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	-13	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	-12	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	-11	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	-10	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	-9	SER	-	EXPRESSION TAG	UNP P0A6L4
B	-8	SER	-	EXPRESSION TAG	UNP P0A6L4
B	-7	GLY	-	EXPRESSION TAG	UNP P0A6L4
B	-6	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	-5	ILE	-	EXPRESSION TAG	UNP P0A6L4
B	-4	GLU	-	EXPRESSION TAG	UNP P0A6L4
B	-3	GLY	-	EXPRESSION TAG	UNP P0A6L4
B	-2	ARG	-	EXPRESSION TAG	UNP P0A6L4
B	-1	HIS	-	EXPRESSION TAG	UNP P0A6L4
B	0	MET	-	EXPRESSION TAG	UNP P0A6L4
B	60	ALA	GLU	ENGINEERED MUTATION	UNP P0A6L4
B	98	HIS	TYR	ENGINEERED MUTATION	UNP P0A6L4
B	115	LEU	PHE	ENGINEERED MUTATION	UNP P0A6L4
B	150	GLY	ASP	ENGINEERED MUTATION	UNP P0A6L4
B	153	TYR	ASN	ENGINEERED MUTATION	UNP P0A6L4
B	251	ILE	VAL	ENGINEERED MUTATION	UNP P0A6L4
B	265	ILE	VAL	ENGINEERED MUTATION	UNP P0A6L4
B	281	CYS	TYR	ENGINEERED MUTATION	UNP P0A6L4
C	-21	MET	-	EXPRESSION TAG	UNP P0A6L4
C	-20	GLY	-	EXPRESSION TAG	UNP P0A6L4
C	-19	HIS	-	EXPRESSION TAG	UNP P0A6L4

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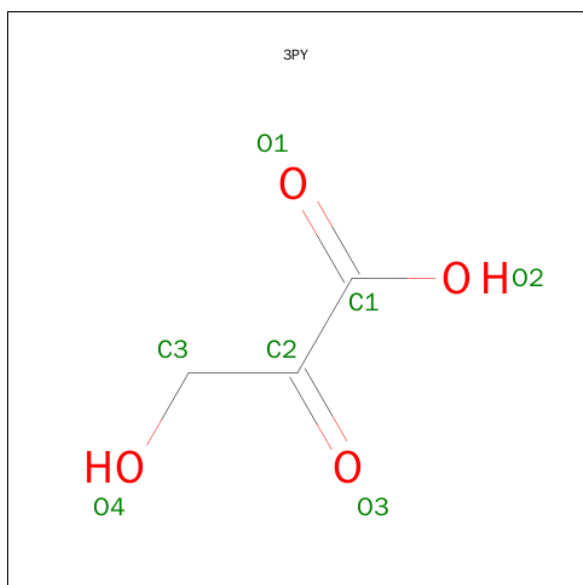
Chain	Residue	Modelled	Actual	Comment	Reference
C	-18	HIS	-	EXPRESSION TAG	UNP P0A6L4
C	-17	HIS	-	EXPRESSION TAG	UNP P0A6L4
C	-16	HIS	-	EXPRESSION TAG	UNP P0A6L4
C	-15	HIS	-	EXPRESSION TAG	UNP P0A6L4
C	-14	HIS	-	EXPRESSION TAG	UNP P0A6L4
C	-13	HIS	-	EXPRESSION TAG	UNP P0A6L4
C	-12	HIS	-	EXPRESSION TAG	UNP P0A6L4
C	-11	HIS	-	EXPRESSION TAG	UNP P0A6L4
C	-10	HIS	-	EXPRESSION TAG	UNP P0A6L4
C	-9	SER	-	EXPRESSION TAG	UNP P0A6L4
C	-8	SER	-	EXPRESSION TAG	UNP P0A6L4
C	-7	GLY	-	EXPRESSION TAG	UNP P0A6L4
C	-6	HIS	-	EXPRESSION TAG	UNP P0A6L4
C	-5	ILE	-	EXPRESSION TAG	UNP P0A6L4
C	-4	GLU	-	EXPRESSION TAG	UNP P0A6L4
C	-3	GLY	-	EXPRESSION TAG	UNP P0A6L4
C	-2	ARG	-	EXPRESSION TAG	UNP P0A6L4
C	-1	HIS	-	EXPRESSION TAG	UNP P0A6L4
C	0	MET	-	EXPRESSION TAG	UNP P0A6L4
C	60	ALA	GLU	ENGINEERED MUTATION	UNP P0A6L4
C	98	HIS	TYR	ENGINEERED MUTATION	UNP P0A6L4
C	115	LEU	PHE	ENGINEERED MUTATION	UNP P0A6L4
C	150	GLY	ASP	ENGINEERED MUTATION	UNP P0A6L4
C	153	TYR	ASN	ENGINEERED MUTATION	UNP P0A6L4
C	251	ILE	VAL	ENGINEERED MUTATION	UNP P0A6L4
C	265	ILE	VAL	ENGINEERED MUTATION	UNP P0A6L4
C	281	CYS	TYR	ENGINEERED MUTATION	UNP P0A6L4
D	-21	MET	-	EXPRESSION TAG	UNP P0A6L4
D	-20	GLY	-	EXPRESSION TAG	UNP P0A6L4
D	-19	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	-18	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	-17	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	-16	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	-15	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	-14	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	-13	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	-12	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	-11	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	-10	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	-9	SER	-	EXPRESSION TAG	UNP P0A6L4
D	-8	SER	-	EXPRESSION TAG	UNP P0A6L4
D	-7	GLY	-	EXPRESSION TAG	UNP P0A6L4

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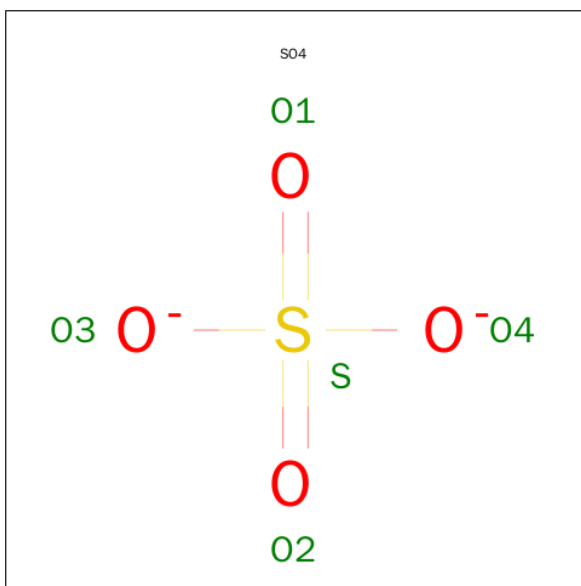
Chain	Residue	Modelled	Actual	Comment	Reference
D	-6	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	-5	ILE	-	EXPRESSION TAG	UNP P0A6L4
D	-4	GLU	-	EXPRESSION TAG	UNP P0A6L4
D	-3	GLY	-	EXPRESSION TAG	UNP P0A6L4
D	-2	ARG	-	EXPRESSION TAG	UNP P0A6L4
D	-1	HIS	-	EXPRESSION TAG	UNP P0A6L4
D	0	MET	-	EXPRESSION TAG	UNP P0A6L4
D	60	ALA	GLU	ENGINEERED MUTATION	UNP P0A6L4
D	98	HIS	TYR	ENGINEERED MUTATION	UNP P0A6L4
D	115	LEU	PHE	ENGINEERED MUTATION	UNP P0A6L4
D	150	GLY	ASP	ENGINEERED MUTATION	UNP P0A6L4
D	153	TYR	ASN	ENGINEERED MUTATION	UNP P0A6L4
D	251	ILE	VAL	ENGINEERED MUTATION	UNP P0A6L4
D	265	ILE	VAL	ENGINEERED MUTATION	UNP P0A6L4
D	281	CYS	TYR	ENGINEERED MUTATION	UNP P0A6L4

- Molecule 2 is 3-HYDROXYPYRUVIC ACID (three-letter code: 3PY) (formula: $C_3H_4O_4$).



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

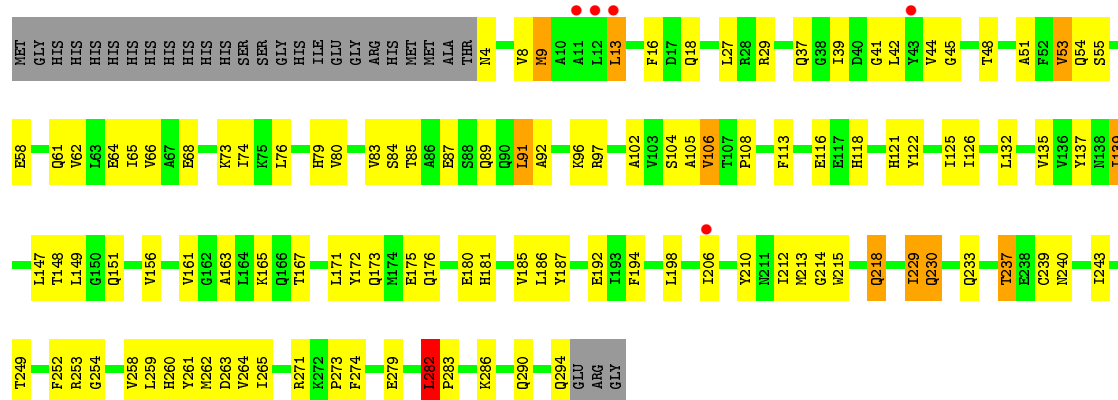
- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	O	S	0	0
			5	4	1		
3	B	1	Total	O	S	0	0
			5	4	1		
3	C	1	Total	O	S	0	0
			5	4	1		
3	D	1	Total	O	S	0	0
			5	4	1		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	146	Total	O	0	0
			146	146		
4	B	113	Total	O	0	0
			113	113		
4	C	108	Total	O	0	0
			108	108		
4	D	126	Total	O	0	0
			126	126		



4 Data and refinement statistics

Property	Value	Source
Space group	P 32 2 1	Depositor
Cell constants a, b, c, α , β , γ	120.83Å 120.83Å 195.88Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	27.83 – 2.35 27.83 – 2.35	Depositor EDS
% Data completeness (in resolution range)	90.6 (27.83-2.35) 95.4 (27.83-2.35)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.41 (at 2.36Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.236 , 0.274 0.227 , 0.267	Depositor DCC
R_{free} test set	3366 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	43.0	Xtriage
Anisotropy	0.037	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 52.9	EDS
Estimated twinning fraction	0.024 for -h,-k,l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 66356 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	9473	wwPDB-VP
Average B, all atoms (Å ²)	44.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.95% 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: 3PY, SO4

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	0.37	0/2273	0.63	1/3076 (0.0%)
1	B	0.37	1/2273 (0.0%)	0.62	1/3076 (0.0%)
1	C	0.35	0/2273	0.63	1/3076 (0.0%)
1	D	0.40	1/2273 (0.0%)	0.64	1/3076 (0.0%)
All	All	0.37	2/9092 (0.0%)	0.63	4/12304 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	282	LEU	C-N	8.08	1.49	1.34
1	B	282	LEU	C-N	7.05	1.47	1.34

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	13	LEU	CA-CB-CG	6.67	130.63	115.30
1	D	13	LEU	CA-CB-CG	6.43	130.09	115.30
1	B	13	LEU	CA-CB-CG	5.62	128.24	115.30
1	C	281	CYS	O-C-N	-5.14	114.48	122.70

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	2234	0	2268	103	0
1	B	2234	0	2268	77	0
1	C	2234	0	2268	100	0
1	D	2234	0	2268	102	0
2	A	6	0	2	0	0
2	B	6	0	2	0	0
2	C	6	0	2	0	0
2	D	6	0	2	0	0
3	A	5	0	0	0	0
3	B	5	0	0	0	0
3	C	5	0	0	1	0
3	D	5	0	0	1	0
4	A	146	0	0	3	0
4	B	113	0	0	1	0
4	C	108	0	0	3	0
4	D	126	0	0	9	0
All	All	9473	0	9080	365	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 21.

All (365) 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:C:80:VAL:HG13	1:C:91:LEU:HB3	1.29	1.10
1:A:218:GLN:HE21	1:A:218:GLN:HA	1.29	0.95
1:A:20:GLN:HE22	1:A:272:LYS:H	1.05	0.94
1:A:80:VAL:HG13	1:A:91:LEU:HB3	1.51	0.91
1:D:192:GLU:HG3	1:D:243:ILE:HD13	1.56	0.87
1:C:218:GLN:HA	1:C:218:GLN:HE21	1.38	0.86
1:A:213:MET:HE1	1:A:216:ARG:HH11	1.41	0.86
1:B:80:VAL:HG13	1:B:91:LEU:CB	2.08	0.83
1:B:115:LEU:HD11	1:B:154:THR:HG21	1.61	0.82
1:A:213:MET:HE3	1:A:216:ARG:HD2	1.63	0.81
1:C:80:VAL:HG13	1:C:91:LEU:CB	2.12	0.80
1:B:218:GLN:HE21	1:B:218:GLN:HA	1.45	0.80
1:C:79:HIS:HE1	1:C:106:VAL:H	1.31	0.79
1:C:56:LEU:HD11	1:C:87:GLU:HG2	1.64	0.79
1:D:106:VAL:HG13	1:D:137:TYR:CD2	2.17	0.79
1:B:80:VAL:HG13	1:B:91:LEU:HB3	1.65	0.78
1:D:108:PRO:HD2	1:D:118:HIS:HD2	1.47	0.78
1:D:258:VAL:O	1:D:262:MET:HG3	1.86	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:PRO:HD2	1:A:118:HIS:HD2	1.51	0.75
1:C:106:VAL:HG13	1:C:137:TYR:CD2	2.20	0.75
1:A:70:ALA:HB1	1:A:74:ILE:CD1	2.16	0.75
1:D:249:THR:O	1:D:253:ARG:HD2	1.88	0.74
1:C:279:GLU:HA	1:C:282:LEU:HD23	1.67	0.74
1:A:121:HIS:O	1:A:125:ILE:HG12	1.87	0.74
1:D:79:HIS:HE1	1:D:106:VAL:H	1.36	0.74
1:A:20:GLN:NE2	1:A:272:LYS:H	1.84	0.74
1:B:51:ALA:HA	1:B:54:GLN:HE21	1.53	0.73
1:A:249:THR:O	1:A:253:ARG:HD2	1.89	0.72
1:C:218:GLN:HA	1:C:218:GLN:NE2	2.04	0.72
1:A:80:VAL:HG13	1:A:91:LEU:CB	2.20	0.71
1:A:108:PRO:HD2	1:A:118:HIS:CD2	2.25	0.71
1:D:192:GLU:HA	1:D:240:ASN:HD21	1.55	0.71
1:C:271:ARG:O	1:C:274:PHE:HB2	1.91	0.70
1:D:279:GLU:HA	1:D:282:LEU:CD2	2.20	0.70
1:C:261:TYR:CZ	1:C:286:LYS:HG2	2.28	0.69
1:C:79:HIS:CE1	1:C:106:VAL:HG22	2.27	0.69
1:A:79:HIS:HE1	1:A:106:VAL:H	1.41	0.69
1:D:233:GLN:O	1:D:237:THR:HG23	1.93	0.69
1:D:79:HIS:CE1	1:D:106:VAL:HG22	2.27	0.68
1:D:264:VAL:HG12	1:D:265:ILE:HD12	1.76	0.68
1:A:186:LEU:H	1:A:203:ASP:HB2	1.58	0.68
1:B:62:VAL:O	1:B:66:VAL:HG12	1.94	0.68
1:C:129:ALA:O	1:C:132:LEU:HD13	1.94	0.67
1:A:114:SER:OG	1:A:117:GLU:HG3	1.95	0.67
1:B:9:MET:O	1:B:206:ILE:HA	1.93	0.67
1:C:50:GLU:O	1:C:53:VAL:HG12	1.95	0.67
1:A:20:GLN:HE22	1:A:272:LYS:N	1.88	0.66
1:A:192:GLU:HG3	1:A:243:ILE:HG21	1.75	0.66
1:A:20:GLN:NE2	1:A:271:ARG:HA	2.10	0.66
1:D:163:ALA:HB2	1:D:185:VAL:HB	1.76	0.66
1:A:218:GLN:NE2	1:A:218:GLN:HA	2.06	0.66
1:B:80:VAL:HG13	1:B:91:LEU:HB2	1.78	0.66
1:C:286:LYS:O	1:C:290:GLN:HG2	1.96	0.65
1:A:192:GLU:HA	1:A:240:ASN:HD21	1.59	0.65
1:D:192:GLU:HA	1:D:240:ASN:ND2	2.11	0.65
1:A:213:MET:HE1	1:A:216:ARG:NH1	2.12	0.65
1:B:192:GLU:HG3	1:B:243:ILE:HG21	1.79	0.65
1:C:79:HIS:HE1	1:C:106:VAL:N	1.94	0.65
1:C:282:LEU:O	1:C:286:LYS:HG3	1.96	0.64

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:149:LEU:HD21	1:C:180:GLU:OE2	1.98	0.64
1:D:106:VAL:HG13	1:D:137:TYR:CE2	2.32	0.64
1:B:242:VAL:HG13	1:B:288:LEU:HD21	1.80	0.64
1:D:279:GLU:HA	1:D:282:LEU:HD22	1.80	0.63
1:D:237:THR:HG22	4:D:457:HOH:O	1.98	0.63
1:A:20:GLN:HE21	1:A:271:ARG:HA	1.63	0.63
1:B:148:THR:O	1:B:152:ILE:HG12	1.99	0.63
1:A:80:VAL:CG1	1:A:88:SER:O	2.47	0.63
1:D:108:PRO:HD2	1:D:118:HIS:CD2	2.32	0.63
1:B:279:GLU:HA	1:B:282:LEU:CD2	2.29	0.63
1:C:192:GLU:HG3	1:C:243:ILE:HD13	1.81	0.62
1:C:152:ILE:O	1:C:156:VAL:HG22	1.99	0.62
1:A:213:MET:HE2	1:A:239:CYS:HA	1.80	0.62
1:D:116:GLU:HG3	4:D:771:HOH:O	1.99	0.62
1:B:236:GLN:HE21	1:B:240:ASN:HD21	1.47	0.62
1:C:149:LEU:HD12	1:C:176:GLN:NE2	2.14	0.62
1:D:156:VAL:HA	1:D:161:VAL:HG11	1.80	0.62
1:D:192:GLU:HG3	1:D:243:ILE:HG21	1.82	0.62
1:C:218:GLN:HE21	1:C:218:GLN:CA	2.05	0.61
1:C:137:TYR:CD1	1:C:165:LYS:HD3	2.35	0.61
1:C:249:THR:O	1:C:253:ARG:HD2	1.99	0.61
1:C:243:ILE:O	1:C:247:ILE:HG12	2.00	0.61
1:A:13:LEU:HD13	1:A:48:THR:HG22	1.83	0.61
1:B:171:LEU:HD12	1:D:171:LEU:HD12	1.82	0.61
1:B:83:VAL:HG23	4:B:402:HOH:O	2.00	0.61
1:D:259:LEU:HB3	1:D:265:ILE:HD13	1.83	0.61
1:D:279:GLU:O	1:D:282:LEU:HD22	2.01	0.61
1:D:173:GLN:HA	1:D:176:GLN:HE21	1.66	0.60
1:A:192:GLU:CG	1:A:243:ILE:HG21	2.31	0.60
1:C:192:GLU:HG3	1:C:243:ILE:HG21	1.83	0.60
1:C:167:THR:HA	1:C:189:GLY:HA3	1.84	0.60
1:B:106:VAL:HA	1:B:137:TYR:HB3	1.83	0.60
1:B:236:GLN:HE21	1:B:240:ASN:ND2	2.00	0.59
1:C:229:ILE:HG13	1:C:230:GLN:N	2.18	0.59
1:A:148:THR:OG1	1:A:151:GLN:HG3	2.02	0.59
1:D:139:ILE:HD13	1:D:167:THR:CG2	2.33	0.59
1:C:236:GLN:HE21	1:C:240:ASN:HD21	1.50	0.59
1:D:229:ILE:HD13	1:D:229:ILE:N	2.17	0.59
1:C:5:LEU:HD22	1:C:5:LEU:H	1.67	0.58
1:D:213:MET:HE1	1:D:239:CYS:HA	1.86	0.58
1:A:163:ALA:HB2	1:A:185:VAL:HB	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:175:GLU:HG2	1:C:240:ASN:OD1	2.03	0.58
1:A:137:TYR:CD1	1:A:165:LYS:HD3	2.39	0.58
1:B:76:LEU:HB3	1:B:100:PHE:CD2	2.38	0.58
1:C:236:GLN:HE21	1:C:240:ASN:ND2	2.02	0.57
1:A:70:ALA:HB1	1:A:74:ILE:HD13	1.85	0.57
1:D:29:ARG:HD2	1:D:264:VAL:O	2.03	0.57
1:D:62:VAL:O	1:D:66:VAL:HG12	2.05	0.57
1:D:218:GLN:HA	1:D:218:GLN:HE21	1.69	0.57
1:A:79:HIS:HE1	1:A:106:VAL:N	2.03	0.57
1:C:80:VAL:CG1	1:C:88:SER:O	2.53	0.57
1:C:149:LEU:O	1:C:149:LEU:HD23	2.05	0.57
1:B:5:LEU:H	1:B:5:LEU:HD23	1.69	0.57
1:A:253:ARG:HD3	1:A:277:VAL:HG22	1.87	0.56
1:A:56:LEU:HD21	1:A:87:GLU:HB3	1.87	0.56
1:D:41:GLY:HA2	1:D:74:ILE:HB	1.86	0.56
1:B:113:PHE:CZ	1:C:273:PRO:HG2	2.40	0.56
1:C:288:LEU:O	1:C:292:LEU:HD23	2.05	0.56
1:C:294:GLN:HG3	1:C:294:GLN:O	2.05	0.56
1:D:61:GLN:O	1:D:65:ILE:HG12	2.05	0.56
1:B:259:LEU:HB3	1:B:265:ILE:HG12	1.88	0.56
1:B:108:PRO:HD2	1:B:118:HIS:HD2	1.70	0.56
1:A:192:GLU:HA	1:A:240:ASN:ND2	2.20	0.56
1:C:64:GLU:O	1:C:68:GLU:HG3	2.05	0.56
1:A:70:ALA:HB1	1:A:74:ILE:HD11	1.88	0.56
1:B:218:GLN:NE2	1:B:218:GLN:HA	2.17	0.56
1:C:79:HIS:CE1	1:C:106:VAL:H	2.18	0.56
1:B:9:MET:HG3	1:B:41:GLY:C	2.26	0.55
1:A:213:MET:CE	1:A:239:CYS:HA	2.36	0.55
1:B:262:MET:O	1:B:263:ASP:HB2	2.05	0.55
1:B:139:ILE:O	1:B:139:ILE:HG23	2.06	0.55
1:D:212:ILE:HG13	1:D:213:MET:HG2	1.89	0.55
1:A:9:MET:O	1:A:206:ILE:HA	2.06	0.55
1:A:113:PHE:CZ	1:D:273:PRO:HG2	2.42	0.55
1:D:9:MET:HG3	1:D:41:GLY:C	2.27	0.54
1:C:20:GLN:HE22	1:C:272:LYS:HG2	1.72	0.54
1:D:121:HIS:CE1	1:D:125:ILE:HD11	2.42	0.54
1:A:88:SER:HB3	1:A:125:ILE:HD12	1.88	0.54
1:A:271:ARG:O	1:A:274:PHE:HB2	2.07	0.54
1:D:121:HIS:HE1	4:D:544:HOH:O	1.91	0.54
1:A:293:MET:O	1:A:294:GLN:HB2	2.06	0.54
1:D:48:THR:HA	1:D:252:PHE:CE1	2.43	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:79:HIS:HE1	1:D:106:VAL:N	2.05	0.54
1:A:273:PRO:HG2	1:D:113:PHE:CZ	2.43	0.54
1:C:4:ASN:HA	4:C:631:HOH:O	2.06	0.53
1:D:192:GLU:CG	1:D:243:ILE:HG21	2.38	0.53
1:C:137:TYR:CE1	1:C:165:LYS:HD3	2.44	0.53
1:B:13:LEU:HD13	1:B:48:THR:HG22	1.90	0.53
1:B:271:ARG:O	1:B:274:PHE:HB2	2.09	0.53
1:C:80:VAL:HG12	1:C:80:VAL:O	2.07	0.53
1:B:80:VAL:HG11	1:B:88:SER:O	2.08	0.53
1:A:171:LEU:HD12	1:C:171:LEU:HD12	1.90	0.53
1:B:121:HIS:O	1:B:125:ILE:HG13	2.09	0.53
1:B:163:ALA:HB2	1:B:185:VAL:HB	1.90	0.53
1:A:117:GLU:OE2	1:D:253:ARG:NH2	2.42	0.52
1:C:5:LEU:HD22	1:C:5:LEU:N	2.25	0.52
1:D:180:GLU:HG3	4:D:376:HOH:O	2.09	0.52
1:A:139:ILE:HD11	1:A:142:LEU:HD22	1.92	0.52
1:A:26:SER:HA	1:A:29:ARG:NH1	2.24	0.52
1:A:85:THR:HA	1:A:121:HIS:CE1	2.45	0.52
1:D:163:ALA:CB	1:D:185:VAL:HB	2.40	0.52
1:C:192:GLU:H	1:C:192:GLU:CD	2.12	0.52
1:C:148:THR:OG1	1:C:151:GLN:HG3	2.10	0.52
1:D:84:SER:OG	1:D:87:GLU:HG3	2.09	0.52
1:D:192:GLU:CD	1:D:192:GLU:H	2.13	0.52
1:B:35:ILE:HG12	1:B:74:ILE:HD13	1.91	0.52
1:A:113:PHE:CE1	1:D:273:PRO:HG2	2.44	0.52
1:B:178:ARG:NH2	1:B:203:ASP:OD2	2.43	0.52
1:A:217:TYR:O	1:A:221:VAL:HG23	2.09	0.51
1:B:213:MET:HA	1:B:215:TRP:CZ3	2.44	0.51
1:C:155:LEU:O	1:C:158:LEU:HB2	2.11	0.51
1:C:135:VAL:HG22	1:C:163:ALA:HB3	1.92	0.51
1:A:237:THR:HG22	1:A:241:LYS:HE3	1.93	0.51
1:C:56:LEU:HD11	1:C:87:GLU:CG	2.38	0.51
1:A:79:HIS:CE1	1:A:106:VAL:H	2.27	0.51
1:C:173:GLN:HA	1:C:176:GLN:HE21	1.75	0.51
1:A:13:LEU:HD13	1:A:48:THR:CG2	2.40	0.51
1:D:229:ILE:HD13	1:D:229:ILE:H	1.76	0.51
1:D:85:THR:OG1	1:D:121:HIS:HD2	1.93	0.51
1:A:137:TYR:CE1	1:A:165:LYS:HD3	2.46	0.50
1:B:108:PRO:HD2	1:B:118:HIS:CD2	2.47	0.50
1:C:139:ILE:HG23	1:C:139:ILE:O	2.11	0.50
1:D:37:GLN:OE1	1:D:214:GLY:N	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:192:GLU:HG3	1:A:243:ILE:HD13	1.93	0.50
1:C:149:LEU:C	1:C:149:LEU:HD23	2.30	0.50
1:B:294:GLN:O	1:B:294:GLN:HG3	2.11	0.50
1:C:30:LEU:HD13	1:C:259:LEU:HD13	1.93	0.50
1:B:80:VAL:CG1	1:B:88:SER:O	2.59	0.50
1:D:165:LYS:HA	1:D:187:TYR:HB2	1.94	0.50
1:B:55:SER:OG	1:B:58:GLU:HG3	2.11	0.50
1:A:103:VAL:HG12	1:A:104:SER:N	2.26	0.50
1:C:33:PHE:O	1:C:37:GLN:HG3	2.12	0.50
1:C:221:VAL:HG13	4:C:496:HOH:O	2.10	0.50
1:C:119:CYS:HB3	1:C:123:ARG:NH1	2.27	0.50
1:B:249:THR:O	1:B:253:ARG:HD2	2.11	0.49
1:C:192:GLU:CG	1:C:243:ILE:HG21	2.42	0.49
1:A:156:VAL:HG11	1:A:184:LEU:HD22	1.93	0.49
1:B:212:ILE:HD12	1:B:292:LEU:HD11	1.95	0.49
1:C:9:MET:O	1:C:206:ILE:HA	2.12	0.49
1:B:113:PHE:CE1	1:C:273:PRO:HG2	2.47	0.49
1:B:267:VAL:CG1	1:B:269:LEU:HD23	2.42	0.49
1:A:213:MET:CE	1:A:216:ARG:HD2	2.39	0.49
1:C:131:GLY:O	1:C:132:LEU:O	2.31	0.49
1:C:79:HIS:CE1	1:C:105:ALA:HA	2.48	0.49
1:D:37:GLN:HG3	1:D:215:TRP:HE3	1.78	0.49
1:D:8:VAL:HG13	1:D:210:TYR:CE2	2.48	0.49
1:B:192:GLU:HB3	1:D:172:TYR:CD1	2.48	0.48
1:D:96:LYS:HD3	1:D:132:LEU:HD13	1.94	0.48
1:C:80:VAL:HG11	1:C:88:SER:O	2.14	0.48
1:A:80:VAL:HG12	1:A:80:VAL:O	2.13	0.48
1:A:51:ALA:HA	1:A:54:GLN:HE21	1.76	0.48
1:A:163:ALA:CB	1:A:185:VAL:HB	2.42	0.48
1:C:132:LEU:HD12	1:C:132:LEU:N	2.29	0.48
1:C:253:ARG:HG3	1:C:254:GLY:N	2.27	0.48
1:B:236:GLN:NE2	1:B:240:ASN:HD21	2.09	0.48
1:A:198:LEU:HA	1:A:202:ALA:HB3	1.95	0.47
1:D:53:VAL:HG13	1:D:53:VAL:O	2.13	0.47
1:B:175:GLU:HG2	1:D:240:ASN:OD1	2.15	0.47
1:A:9:MET:HB2	1:A:206:ILE:HG12	1.95	0.47
1:B:261:TYR:CD2	1:B:286:LYS:HA	2.49	0.47
1:B:229:ILE:C	1:B:229:ILE:HD13	2.35	0.47
1:B:117:GLU:OE2	1:C:253:ARG:NH2	2.45	0.47
1:B:80:VAL:HG12	1:B:80:VAL:O	2.14	0.47
1:B:51:ALA:HA	1:B:54:GLN:NE2	2.25	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:236:GLN:NE2	1:C:240:ASN:HD21	2.13	0.47
1:C:240:ASN:N	1:C:240:ASN:HD22	2.13	0.47
1:D:13:LEU:HD13	1:D:48:THR:CG2	2.45	0.47
1:D:147:LEU:HA	1:D:151:GLN:OE1	2.14	0.47
1:D:261:TYR:CD1	1:D:286:LYS:HE3	2.49	0.47
1:B:165:LYS:HA	1:B:187:TYR:HB2	1.95	0.47
1:D:80:VAL:HB	1:D:91:LEU:HB3	1.96	0.47
1:B:13:LEU:HD13	1:B:48:THR:CG2	2.44	0.47
1:C:289:ALA:O	1:C:293:MET:HG3	2.15	0.47
1:A:78:ALA:HB2	1:A:100:PHE:CD2	2.49	0.47
1:A:45:GLY:O	1:A:79:HIS:HD2	1.97	0.47
1:B:240:ASN:N	1:B:240:ASN:HD22	2.12	0.47
1:D:85:THR:O	1:D:89:GLN:HG3	2.14	0.47
1:D:51:ALA:HA	1:D:54:GLN:HE21	1.80	0.47
1:B:267:VAL:HG12	1:B:269:LEU:HD23	1.97	0.47
1:A:122:TYR:O	1:A:126:ILE:HG13	2.14	0.47
1:A:80:VAL:HG11	1:A:88:SER:O	2.16	0.46
1:C:256:LYS:HB3	1:C:268:PRO:HA	1.97	0.46
1:D:79:HIS:HE1	1:D:106:VAL:HG22	1.79	0.46
1:D:9:MET:O	1:D:206:ILE:HA	2.16	0.46
1:A:174:MET:SD	1:A:186:LEU:HD13	2.56	0.46
1:D:229:ILE:H	1:D:229:ILE:CD1	2.29	0.46
1:D:18:GLN:HA	1:D:18:GLN:NE2	2.31	0.46
1:D:64:GLU:O	1:D:68:GLU:HG3	2.16	0.46
1:D:122:TYR:O	1:D:126:ILE:HG13	2.16	0.46
1:A:240:ASN:OD1	1:C:175:GLU:HG2	2.16	0.46
1:A:62:VAL:O	1:A:66:VAL:HG23	2.16	0.46
1:A:192:GLU:HB3	1:C:172:TYR:CD1	2.50	0.46
1:B:114:SER:OG	1:B:117:GLU:HG3	2.16	0.45
1:B:59:ARG:O	1:B:63:LEU:HG	2.16	0.45
1:B:70:ALA:O	1:B:71:LYS:C	2.54	0.45
1:D:192:GLU:HG3	1:D:243:ILE:CD1	2.38	0.45
1:A:106:VAL:HA	1:A:137:TYR:HB3	1.98	0.45
1:A:165:LYS:HA	1:A:187:TYR:HB2	1.97	0.45
1:B:292:LEU:O	1:B:294:GLN:N	2.44	0.45
1:A:54:GLN:HA	1:A:271:ARG:HH21	1.81	0.45
1:A:22:LEU:HD22	4:A:445:HOH:O	2.15	0.45
1:A:139:ILE:O	1:A:139:ILE:HG23	2.15	0.45
1:C:258:VAL:HG21	1:C:288:LEU:HD23	1.98	0.45
1:B:121:HIS:CE1	1:B:125:ILE:HD11	2.51	0.45
1:A:28:ARG:HD3	1:A:69:GLU:OE2	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:230:GLN:CA	1:D:230:GLN:HE21	2.28	0.45
1:C:76:LEU:HB3	1:C:100:PHE:CD2	2.51	0.45
1:B:148:THR:HG23	1:B:151:GLN:OE1	2.16	0.45
1:A:171:LEU:HA	1:A:171:LEU:HD23	1.79	0.45
1:B:229:ILE:O	1:B:229:ILE:HD13	2.16	0.45
1:D:55:SER:OG	1:D:58:GLU:HG3	2.17	0.45
1:B:20:GLN:NE2	1:B:271:ARG:HA	2.32	0.45
1:C:14:THR:HG23	1:C:49:GLY:O	2.16	0.44
1:C:52:PHE:HZ	1:C:106:VAL:HG21	1.82	0.44
1:C:79:HIS:NE2	1:C:106:VAL:HG22	2.32	0.44
1:D:253:ARG:HG3	1:D:254:GLY:N	2.33	0.44
1:A:45:GLY:O	1:A:79:HIS:CD2	2.70	0.44
1:D:42:LEU:HG	1:D:74:ILE:HD12	2.00	0.44
1:B:31:VAL:O	1:B:35:ILE:HG13	2.17	0.44
1:A:191:ASP:HB3	1:A:217:TYR:OH	2.16	0.44
1:D:290:GLN:HE21	1:D:290:GLN:HA	1.82	0.44
1:D:16:PHE:CE2	1:D:271:ARG:HG3	2.51	0.44
1:A:196:SER:HB3	1:C:196:SER:HB3	2.00	0.44
1:D:261:TYR:CE1	1:D:286:LYS:HE3	2.52	0.44
1:D:253:ARG:HG2	4:D:551:HOH:O	2.18	0.44
1:C:108:PRO:HD2	1:C:118:HIS:HD2	1.82	0.44
1:C:282:LEU:HD13	1:C:282:LEU:HA	1.58	0.44
1:C:173:GLN:O	1:C:177:ILE:HG13	2.17	0.44
1:C:277:VAL:HG12	1:C:278:ASP:N	2.33	0.44
1:A:59:ARG:O	1:A:63:LEU:HG	2.18	0.44
1:B:90:GLN:HE21	1:B:90:GLN:HB2	1.64	0.44
1:A:80:VAL:HG11	1:A:92:ALA:N	2.33	0.44
1:B:155:LEU:O	1:B:158:LEU:HB2	2.17	0.44
1:B:163:ALA:CB	1:B:185:VAL:HB	2.48	0.44
1:D:8:VAL:HG12	1:D:39:ILE:HD11	2.00	0.44
1:D:271:ARG:O	1:D:274:PHE:HB2	2.18	0.44
1:A:237:THR:HG23	1:C:179:ARG:HD3	1.99	0.43
1:D:104:SER:HB2	1:D:135:VAL:HB	2.00	0.43
1:A:284:GLU:HG3	4:A:769:HOH:O	2.17	0.43
1:D:79:HIS:CE1	1:D:105:ALA:HA	2.54	0.43
1:A:218:GLN:HE21	1:A:218:GLN:CA	2.08	0.43
1:A:249:THR:HG22	1:A:281:CYS:SG	2.59	0.43
1:D:139:ILE:HD12	4:D:675:HOH:O	2.17	0.43
1:A:139:ILE:CD1	1:A:142:LEU:HD22	2.48	0.43
1:A:26:SER:HA	1:A:29:ARG:HH11	1.83	0.43
1:A:153:TYR:HB2	4:A:629:HOH:O	2.19	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:17:ASP:OD1	1:A:19:GLN:N	2.48	0.43
1:D:79:HIS:CE1	1:D:106:VAL:H	2.26	0.43
1:C:261:TYR:OH	1:C:282:LEU:HD12	2.18	0.43
1:C:16:PHE:CE2	1:C:271:ARG:HG3	2.54	0.43
1:D:148:THR:OG1	1:D:151:GLN:HG3	2.18	0.43
1:C:164:LEU:CD2	1:C:166:GLN:HG2	2.48	0.43
1:A:16:PHE:CE2	1:A:271:ARG:HG3	2.54	0.43
1:C:97:ARG:HD2	3:C:647:SO4:O2	2.19	0.43
1:D:282:LEU:N	1:D:283:PRO:HD2	2.34	0.42
1:D:229:ILE:N	1:D:229:ILE:CD1	2.81	0.42
1:B:32:GLN:NE2	1:B:69:GLU:OE1	2.52	0.42
1:C:45:GLY:HA3	4:C:307:HOH:O	2.17	0.42
1:C:288:LEU:HA	1:C:291:GLN:HG2	2.01	0.42
1:A:273:PRO:HG2	1:D:113:PHE:CE1	2.53	0.42
1:B:291:GLN:HG3	1:B:292:LEU:N	2.33	0.42
1:C:103:VAL:CG1	1:C:104:SER:N	2.83	0.42
1:D:233:GLN:O	1:D:237:THR:CG2	2.64	0.42
1:D:102:ALA:HA	1:D:132:LEU:HB3	2.01	0.42
1:A:267:VAL:CG1	1:A:269:LEU:HD23	2.49	0.42
1:D:218:GLN:HA	1:D:218:GLN:NE2	2.33	0.42
1:B:253:ARG:HD3	1:B:277:VAL:HG22	2.02	0.42
1:A:156:VAL:HA	1:A:161:VAL:HG11	2.02	0.42
1:A:88:SER:HB2	1:A:125:ILE:CD1	2.50	0.42
1:B:279:GLU:HA	1:B:282:LEU:HD22	2.02	0.42
1:B:259:LEU:CB	1:B:265:ILE:HG12	2.49	0.42
1:D:181:HIS:HE1	4:D:377:HOH:O	2.02	0.42
1:C:70:ALA:O	1:C:71:LYS:C	2.58	0.42
1:C:106:VAL:HG13	1:C:137:TYR:CE2	2.54	0.42
1:C:51:ALA:HA	1:C:54:GLN:HE21	1.84	0.42
1:C:106:VAL:HA	1:C:137:TYR:HB3	2.01	0.42
1:A:48:THR:HA	1:A:252:PHE:CE1	2.55	0.41
1:C:163:ALA:HB2	1:C:185:VAL:HB	2.02	0.41
1:D:80:VAL:HG11	1:D:92:ALA:HA	2.02	0.41
1:D:18:GLN:HA	1:D:18:GLN:HE21	1.84	0.41
1:B:61:GLN:O	1:B:65:ILE:HD12	2.20	0.41
1:C:8:VAL:O	1:C:40:ASP:HB2	2.21	0.41
1:B:92:ALA:O	1:B:95:ALA:HB3	2.20	0.41
1:D:290:GLN:NE2	1:D:290:GLN:HA	2.35	0.41
1:A:80:VAL:HG12	1:A:88:SER:O	2.20	0.41
1:B:61:GLN:HG2	1:B:65:ILE:CD1	2.50	0.41
1:D:194:PHE:CE2	1:D:198:LEU:HD11	2.55	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:289:ALA:O	1:A:293:MET:HG3	2.20	0.41
1:D:45:GLY:O	1:D:79:HIS:HD2	2.04	0.41
1:B:218:GLN:HE21	1:B:218:GLN:CA	2.18	0.41
1:C:288:LEU:HA	1:C:291:GLN:HE21	1.86	0.41
1:A:291:GLN:O	1:A:294:GLN:N	2.54	0.41
1:B:245:LEU:O	1:B:249:THR:HG23	2.20	0.41
1:D:4:ASN:N	4:D:318:HOH:O	2.54	0.41
1:B:287:ALA:O	1:B:290:GLN:HB2	2.19	0.41
1:D:13:LEU:HD13	1:D:48:THR:HG22	2.02	0.41
1:D:260:HIS:HD2	1:D:265:ILE:O	2.04	0.40
1:A:237:THR:O	1:A:241:LYS:HG3	2.20	0.40
1:C:150:GLY:O	1:C:153:TYR:HB3	2.20	0.40
1:C:249:THR:HG22	1:C:281:CYS:SG	2.61	0.40
1:C:25:ALA:HB1	1:C:29:ARG:HH21	1.86	0.40
1:A:287:ALA:O	1:A:290:GLN:HB2	2.21	0.40
1:D:83:VAL:HG23	4:D:362:HOH:O	2.20	0.40
1:C:253:ARG:HD3	1:C:277:VAL:HA	2.04	0.40
1:A:229:ILE:O	1:A:233:GLN:HG3	2.21	0.40
1:D:97:ARG:HD3	3:D:645:SO4:O2	2.22	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	289/319 (91%)	282 (98%)	7 (2%)	0	100	100
1	B	289/319 (91%)	276 (96%)	13 (4%)	0	100	100
1	C	289/319 (91%)	275 (95%)	12 (4%)	2 (1%)	26	29
1	D	289/319 (91%)	279 (96%)	10 (4%)	0	100	100
All	All	1156/1276 (91%)	1112 (96%)	42 (4%)	2 (0%)	52	63

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	C	132	LEU
1	C	251	ILE

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	238/261 (91%)	227 (95%)	11 (5%)	33	42
1	B	238/261 (91%)	224 (94%)	14 (6%)	24	28
1	C	238/261 (91%)	231 (97%)	7 (3%)	50	64
1	D	238/261 (91%)	219 (92%)	19 (8%)	15	16
All	All	952/1044 (91%)	901 (95%)	51 (5%)	27	33

All (51) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	13	LEU
1	A	27	LEU
1	A	44	VAL
1	A	74	ILE
1	A	76	LEU
1	A	91	LEU
1	A	149	LEU
1	A	175	GLU
1	A	203	ASP
1	A	218	GLN
1	A	234	LYS
1	B	5	LEU
1	B	6	ARG
1	B	9	MET
1	B	91	LEU
1	B	103	VAL
1	B	149	LEU
1	B	158	LEU

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Mol	Chain	Res	Type
1	B	175	GLU
1	B	218	GLN
1	B	229	ILE
1	B	234	LYS
1	B	263	ASP
1	B	271	ARG
1	B	282	LEU
1	C	14	THR
1	C	27	LEU
1	C	106	VAL
1	C	115	LEU
1	C	175	GLU
1	C	218	GLN
1	C	282	LEU
1	D	9	MET
1	D	27	LEU
1	D	44	VAL
1	D	53	VAL
1	D	73	LYS
1	D	76	LEU
1	D	91	LEU
1	D	106	VAL
1	D	139	ILE
1	D	149	LEU
1	D	175	GLU
1	D	186	LEU
1	D	218	GLN
1	D	229	ILE
1	D	230	GLN
1	D	237	THR
1	D	263	ASP
1	D	282	LEU
1	D	294	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (62) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	4	ASN
1	A	18	GLN
1	A	20	GLN
1	A	54	GLN
1	A	61	GLN

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Mol	Chain	Res	Type
1	A	79	HIS
1	A	90	GLN
1	A	118	HIS
1	A	138	ASN
1	A	166	GLN
1	A	218	GLN
1	A	233	GLN
1	A	240	ASN
1	A	260	HIS
1	B	18	GLN
1	B	20	GLN
1	B	37	GLN
1	B	54	GLN
1	B	61	GLN
1	B	90	GLN
1	B	118	HIS
1	B	121	HIS
1	B	138	ASN
1	B	166	GLN
1	B	176	GLN
1	B	218	GLN
1	B	230	GLN
1	B	240	ASN
1	B	260	HIS
1	B	294	GLN
1	C	4	ASN
1	C	18	GLN
1	C	20	GLN
1	C	54	GLN
1	C	61	GLN
1	C	79	HIS
1	C	90	GLN
1	C	118	HIS
1	C	138	ASN
1	C	166	GLN
1	C	176	GLN
1	C	218	GLN
1	C	240	ASN
1	C	260	HIS
1	C	291	GLN
1	C	294	GLN
1	D	18	GLN

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Mol	Chain	Res	Type
1	D	54	GLN
1	D	79	HIS
1	D	90	GLN
1	D	118	HIS
1	D	121	HIS
1	D	138	ASN
1	D	166	GLN
1	D	176	GLN
1	D	181	HIS
1	D	218	GLN
1	D	230	GLN
1	D	240	ASN
1	D	260	HIS
1	D	290	GLN
1	D	294	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no carbohydrates in this entry.

5.6 Ligand geometry [i](#)

8 ligands are modelled in this entry.

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	3PY	A	640	1	2,5,6	2.60	1 (50%)	1,5,7	2.25	1 (100%)
3	SO4	A	644	-	4,4,4	0.12	0	6,6,6	0.16	0
2	3PY	B	641	1	2,5,6	2.51	1 (50%)	1,5,7	2.49	1 (100%)
3	SO4	B	646	-	4,4,4	0.16	0	6,6,6	0.11	0
2	3PY	C	642	1	2,5,6	2.62	1 (50%)	1,5,7	2.24	1 (100%)
3	SO4	C	647	-	4,4,4	0.16	0	6,6,6	0.07	0
2	3PY	D	643	1	2,5,6	2.55	1 (50%)	1,5,7	2.52	1 (100%)
3	SO4	D	645	-	4,4,4	0.14	0	6,6,6	0.21	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	3PY	A	640	1	-	0/1/3/6	0/0/0/0
3	SO4	A	644	-	-	0/0/0/0	0/0/0/0
2	3PY	B	641	1	-	0/1/3/6	0/0/0/0
3	SO4	B	646	-	-	0/0/0/0	0/0/0/0
2	3PY	C	642	1	-	0/1/3/6	0/0/0/0
3	SO4	C	647	-	-	0/0/0/0	0/0/0/0
2	3PY	D	643	1	-	0/1/3/6	0/0/0/0
3	SO4	D	645	-	-	0/0/0/0	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	640	3PY	O4-C3	-3.27	1.24	1.42
2	C	642	3PY	O4-C3	-3.25	1.24	1.42
2	D	643	3PY	O4-C3	-3.22	1.24	1.42
2	B	641	3PY	O4-C3	-3.17	1.24	1.42

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	643	3PY	O4-C3-C2	-2.52	102.95	110.72
2	B	641	3PY	O4-C3-C2	-2.49	103.04	110.72
2	A	640	3PY	O4-C3-C2	-2.25	103.79	110.72
2	C	642	3PY	O4-C3-C2	-2.24	103.81	110.72

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	C	647	SO4	1	0
3	D	645	SO4	1	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	291/319 (91%)	-0.06	5 (1%)	73 83	28, 39, 54, 65	16 (5%)
1	B	291/319 (91%)	0.26	19 (6%)	22 33	34, 44, 63, 76	16 (5%)
1	C	291/319 (91%)	0.29	22 (7%)	17 26	30, 48, 65, 86	16 (5%)
1	D	291/319 (91%)	0.01	5 (1%)	73 83	26, 38, 54, 68	13 (4%)
All	All	1164/1276 (91%)	0.12	51 (4%)	38 52	26, 42, 62, 86	61 (5%)

All (51) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	153	TYR	4.0
1	C	206	ILE	3.7
1	B	11	ALA	3.3
1	C	46	GLY	3.3
1	B	43	TYR	3.2
1	B	48	THR	3.1
1	C	292	LEU	3.1
1	B	227	GLY	3.1
1	B	135	VAL	3.0
1	C	137	TYR	2.8
1	C	105	ALA	2.8
1	C	70	ALA	2.7
1	B	10	ALA	2.6
1	C	48	THR	2.6
1	D	206	ILE	2.6
1	C	106	VAL	2.6
1	B	282	LEU	2.6
1	D	13	LEU	2.6
1	C	104	SER	2.5
1	B	12	LEU	2.5
1	C	72	GLY	2.5

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Mol	Chain	Res	Type	RSRZ
1	D	11	ALA	2.5
1	C	136	VAL	2.4
1	B	72	GLY	2.4
1	C	290	GLN	2.4
1	B	148	THR	2.3
1	C	294	GLN	2.3
1	B	208	SER	2.3
1	C	230	GLN	2.3
1	A	72	GLY	2.3
1	C	80	VAL	2.3
1	C	286	LYS	2.3
1	C	18	GLN	2.3
1	B	263	ASP	2.3
1	C	278	ASP	2.3
1	B	165	LYS	2.2
1	A	11	ALA	2.2
1	C	43	TYR	2.2
1	D	43	TYR	2.2
1	C	207	GLY	2.2
1	B	206	ILE	2.2
1	B	13	LEU	2.2
1	C	79	HIS	2.2
1	B	77	ILE	2.1
1	A	180	GLU	2.1
1	C	47	SER	2.1
1	B	18	GLN	2.0
1	A	10	ALA	2.0
1	D	12	LEU	2.0
1	A	43	TYR	2.0
1	B	207	GLY	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 [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
2	3PY	D	643	6/7	0.88	0.45	5.70	55,56,59,62	0
2	3PY	A	640	6/7	0.92	0.41	4.57	46,50,54,57	0
2	3PY	B	641	6/7	0.93	0.42	2.60	52,54,58,61	0
2	3PY	C	642	6/7	0.95	0.39	2.54	49,50,56,63	0
3	SO4	B	646	5/5	0.91	0.16	0.04	57,57,58,58	5
3	SO4	D	645	5/5	0.99	0.10	-0.36	42,44,45,47	0
3	SO4	A	644	5/5	0.98	0.13	-0.46	42,43,44,46	5
3	SO4	C	647	5/5	0.96	0.14	-	47,49,50,50	5

6.5 Other polymers [i](#)

There are no such residues in this entry.