



Full wwPDB X-ray Structure Validation Report

Feb 27, 2014 – 03:02 AM GMT

PDB ID : 3DTY
Title : Crystal structure of an Oxidoreductase from Pseudomonas syringae
Authors : Eswaramoorthy, S.; Mahmood, A.; Burley, S.K.; Swaminathan, S.; New York
SGX Research Center for Structural Genomics (NYSGXRC)
Deposited on : 2008-07-16
Resolution : 2.04 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>

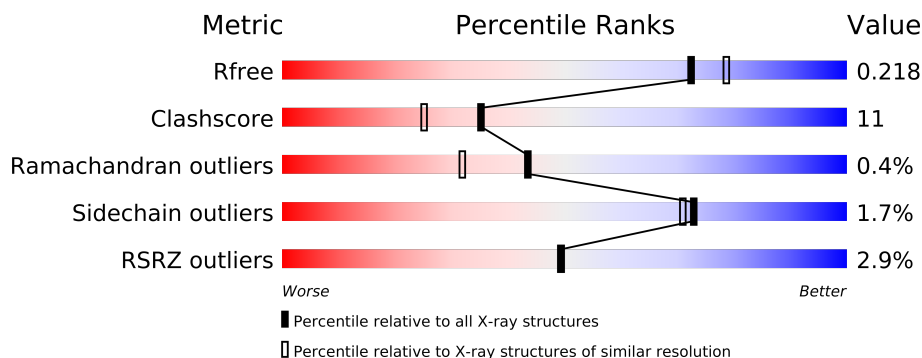
The following versions of software and data (see [references](#)) were used in the production of this report:

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 2.04 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	1380 (2.06-2.02)
Clashscore	79885	1577 (2.06-2.02)
Ramachandran outliers	78287	1565 (2.06-2.02)
Sidechain outliers	78261	1565 (2.06-2.02)
RSRZ outliers	66119	1381 (2.06-2.02)

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	398	
1	B	398	
1	D	398	
1	E	398	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12349 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 Oxidoreductase, Gfo/Idh/MocA family.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	374	Total	C	N	O	S	Se	0	0	0
			2914	1826	530	539	6	13			
1	B	364	Total	C	N	O	S	Se	0	0	0
			2818	1771	502	526	6	13			
1	D	362	Total	C	N	O	S	Se	0	0	0
			2802	1762	498	523	6	13			
1	E	374	Total	C	N	O	S	Se	0	0	0
			2914	1826	530	539	6	13			

There are 44 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	MSE	-	expression tag	UNP Q880Y1
A	1	SER	-	expression tag	UNP Q880Y1
A	2	LEU	-	expression tag	UNP Q880Y1
A	390	GLU	-	expression tag	UNP Q880Y1
A	391	GLY	-	expression tag	UNP Q880Y1
A	392	HIS	-	expression tag	UNP Q880Y1
A	393	HIS	-	expression tag	UNP Q880Y1
A	394	HIS	-	expression tag	UNP Q880Y1
A	395	HIS	-	expression tag	UNP Q880Y1
A	396	HIS	-	expression tag	UNP Q880Y1
A	397	HIS	-	expression tag	UNP Q880Y1
B	0	MSE	-	expression tag	UNP Q880Y1
B	1	SER	-	expression tag	UNP Q880Y1
B	2	LEU	-	expression tag	UNP Q880Y1
B	390	GLU	-	expression tag	UNP Q880Y1
B	391	GLY	-	expression tag	UNP Q880Y1
B	392	HIS	-	expression tag	UNP Q880Y1
B	393	HIS	-	expression tag	UNP Q880Y1
B	394	HIS	-	expression tag	UNP Q880Y1
B	395	HIS	-	expression tag	UNP Q880Y1
B	396	HIS	-	expression tag	UNP Q880Y1

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Chain	Residue	Modelled	Actual	Comment	Reference
B	397	HIS	-	expression tag	UNP Q880Y1
D	0	MSE	-	expression tag	UNP Q880Y1
D	1	SER	-	expression tag	UNP Q880Y1
D	2	LEU	-	expression tag	UNP Q880Y1
D	390	GLU	-	expression tag	UNP Q880Y1
D	391	GLY	-	expression tag	UNP Q880Y1
D	392	HIS	-	expression tag	UNP Q880Y1
D	393	HIS	-	expression tag	UNP Q880Y1
D	394	HIS	-	expression tag	UNP Q880Y1
D	395	HIS	-	expression tag	UNP Q880Y1
D	396	HIS	-	expression tag	UNP Q880Y1
D	397	HIS	-	expression tag	UNP Q880Y1
E	0	MSE	-	expression tag	UNP Q880Y1
E	1	SER	-	expression tag	UNP Q880Y1
E	2	LEU	-	expression tag	UNP Q880Y1
E	390	GLU	-	expression tag	UNP Q880Y1
E	391	GLY	-	expression tag	UNP Q880Y1
E	392	HIS	-	expression tag	UNP Q880Y1
E	393	HIS	-	expression tag	UNP Q880Y1
E	394	HIS	-	expression tag	UNP Q880Y1
E	395	HIS	-	expression tag	UNP Q880Y1
E	396	HIS	-	expression tag	UNP Q880Y1
E	397	HIS	-	expression tag	UNP Q880Y1

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	2	Total Mg 2 2	0	0
2	E	2	Total Mg 2 2	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	248	Total O 248 248	0	0
3	B	188	Total O 188 188	0	0
3	D	198	Total O 198 198	0	0

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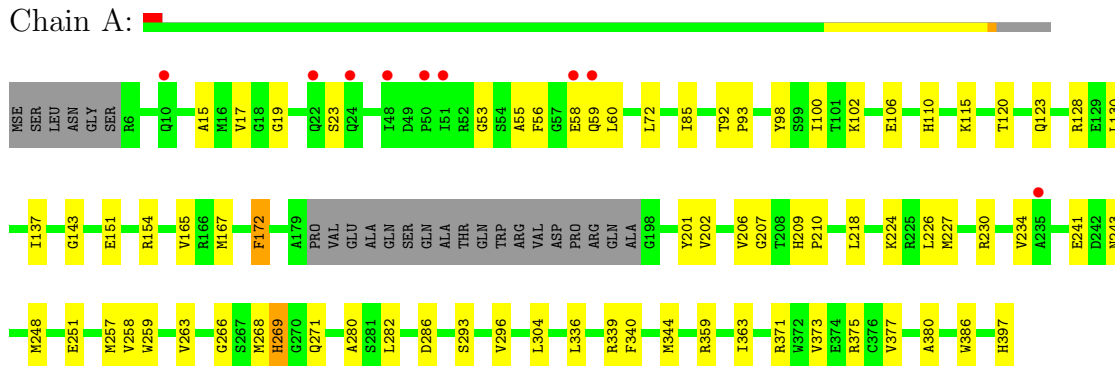
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	263	Total 263	O 263	0	0

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.

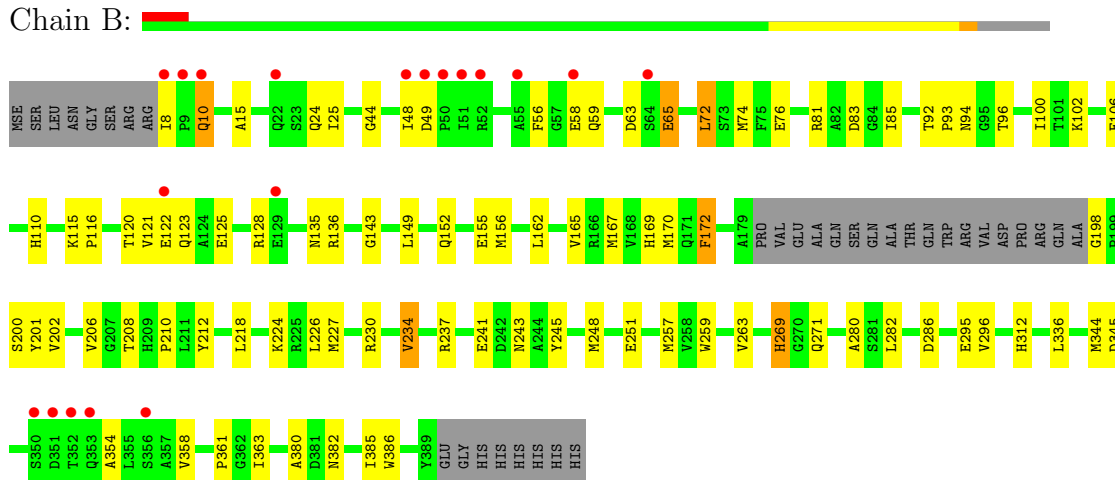
- Molecule 1: Oxidoreductase, Gfo/Idh/MocA family

Chain A:



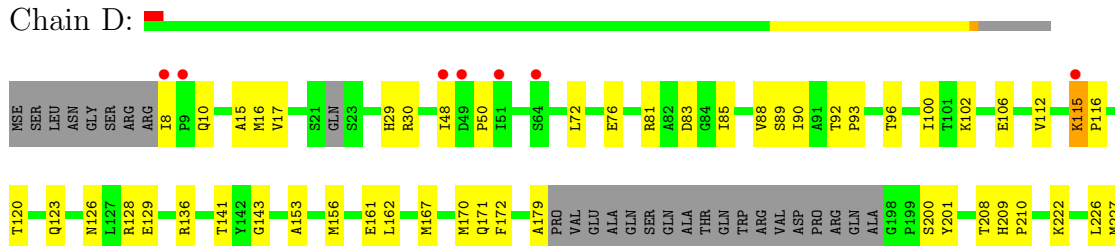
- Molecule 1: Oxidoreductase, Gfo/Idh/MocA family

Chain B:



- Molecule 1: Oxidoreductase, Gfo/Idh/MocA family

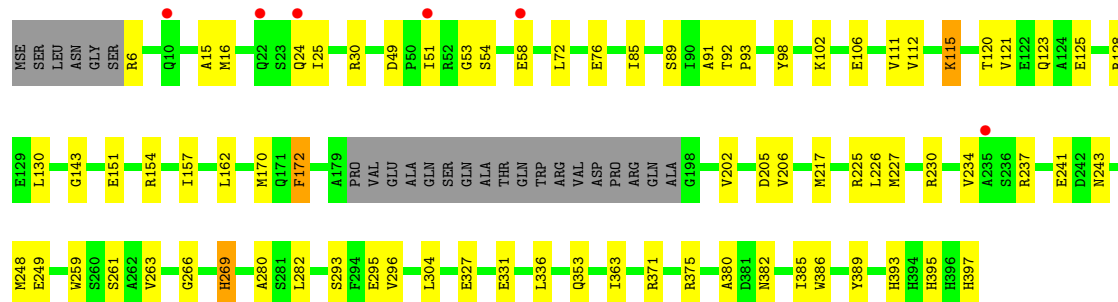
Chain D:





- Molecule 1: Oxidoreductase, Gfo/Idh/MocA family

Chain E:



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	94.28Å 80.93Å 110.54Å 90.00° 93.63° 90.00°	Depositor
Resolution (Å)	50.00 – 2.04 33.48 – 2.03	Depositor EDS
% Data completeness (in resolution range)	94.4 (50.00-2.04) 94.2 (33.48-2.03)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.03Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.193 , 0.218 0.193 , 0.218	Depositor DCC
R_{free} test set	4034 reflections (4.04%)	DCC
Wilson B-factor (Å ²)	14.3	Xtriage
Anisotropy	0.139	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.42 , 35.8	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	2 of 104582 reflections (0.002%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	12349	wwPDB-VP
Average B, all atoms (Å ²)	17.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.04% of the height of the origin peak. No significant pseudotranslation is detected.*

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: MG

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	0.34	0/2972	0.62	1/4002 (0.0%)
1	B	0.32	0/2870	0.61	1/3867 (0.0%)
1	D	0.32	0/2852	0.61	0/3839
1	E	0.33	0/2972	0.62	0/4002
All	All	0.33	0/11666	0.61	2/15710 (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	B	234	VAL	N-CA-C	-5.07	97.32	111.00
1	A	234	VAL	N-CA-C	-5.03	97.43	111.00

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	2914	0	2792	52	0
1	B	2818	0	2715	70	0
1	D	2802	0	2696	74	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	E	2914	0	2792	52	0
2	A	2	0	0	0	0
2	E	2	0	0	0	0
3	A	248	0	0	2	0
3	B	188	0	0	3	0
3	D	198	0	0	9	0
3	E	263	0	0	7	0
All	All	12349	0	10995	237	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 (237) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:115:LYS:HD3	3:D:576:HOH:O	1.49	1.10
1:E:6:ARG:HB3	3:E:575:HOH:O	1.55	1.05
1:E:151:GLU:HG2	3:E:573:HOH:O	1.70	0.92
1:A:243:ASN:HD21	1:A:259:TRP:HE1	1.19	0.89
1:E:243:ASN:HD21	1:E:259:TRP:HE1	1.20	0.89
1:D:16:MSE:HE1	1:D:29:HIS:HB3	1.59	0.84
1:D:8:ILE:HG21	3:D:526:HOH:O	1.77	0.84
1:D:15:ALA:HB2	1:D:85:ILE:HG21	1.57	0.84
1:B:63:ASP:HB3	1:B:65:GLU:OE1	1.78	0.83
1:A:227:MSE:HE3	1:A:386:TRP:CE2	2.14	0.83
1:B:245:TYR:HB3	1:B:257:MSE:HE1	1.62	0.81
1:D:243:ASN:HD21	1:D:259:TRP:HE1	1.28	0.80
1:D:227:MSE:HE3	1:D:386:TRP:CE2	2.16	0.80
1:D:8:ILE:CG2	3:D:526:HOH:O	2.29	0.79
1:B:227:MSE:HE3	1:B:386:TRP:CE2	2.17	0.79
1:B:243:ASN:HD21	1:B:259:TRP:HE1	1.31	0.77
1:A:230:ARG:HG2	1:A:380:ALA:HB1	1.68	0.76
1:D:115:LYS:HZ3	1:D:209:HIS:CE1	2.03	0.75
1:A:154:ARG:CZ	1:A:359:ARG:HD3	2.18	0.73
1:B:122:GLU:HG3	1:B:123:GLN:N	2.04	0.72
1:A:55:ALA:O	1:A:58:GLU:HG2	1.89	0.72
1:B:122:GLU:HG3	1:B:123:GLN:H	1.53	0.72
1:A:110:HIS:HD2	1:A:137:ILE:H	1.37	0.71
1:E:102:LYS:O	1:E:106:GLU:HG3	1.90	0.71
1:E:230:ARG:HG2	1:E:380:ALA:HB1	1.73	0.71
1:A:102:LYS:O	1:A:106:GLU:HG3	1.92	0.70
1:B:162:LEU:HD11	1:B:282:LEU:HG	1.73	0.69
1:E:49:ASP:OD2	1:E:51:ILE:HG13	1.92	0.69

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:E:227:MSE:HE3	1:E:386:TRP:CE2	2.27	0.69
1:D:359:ARG:HA	3:D:568:HOH:O	1.94	0.67
1:A:151:GLU:HG2	3:A:494:HOH:O	1.94	0.67
1:B:121:VAL:O	1:B:125:GLU:HG3	1.96	0.66
1:E:6:ARG:CB	3:E:575:HOH:O	2.27	0.65
1:E:53:GLY:HA3	3:E:649:HOH:O	1.96	0.65
1:E:154:ARG:HD2	3:E:440:HOH:O	1.98	0.64
1:A:15:ALA:HB2	1:A:85:ILE:HG21	1.80	0.64
1:B:115:LYS:HB3	1:B:115:LYS:NZ	2.14	0.62
1:B:120:THR:OG1	1:B:123:GLN:HG3	2.00	0.62
1:D:115:LYS:HE3	1:D:208:THR:OG1	1.99	0.62
1:A:92:THR:HB	1:A:93:PRO:HD2	1.82	0.61
1:D:16:MSE:HE3	1:D:89:SER:HB2	1.83	0.61
1:D:200:SER:HB2	1:D:241:GLU:OE2	2.03	0.59
1:B:15:ALA:HB2	1:B:85:ILE:HG21	1.83	0.59
1:D:72:LEU:O	1:D:76:GLU:HG2	2.03	0.59
1:E:15:ALA:HB2	1:E:85:ILE:HG21	1.84	0.59
1:E:72:LEU:O	1:E:76:GLU:HG3	2.03	0.58
1:B:243:ASN:ND2	1:B:259:TRP:HE1	2.00	0.58
1:E:327:GLU:HG2	1:E:331:GLU:HG2	1.84	0.58
1:D:317:ILE:HD11	3:D:498:HOH:O	2.02	0.58
1:E:385:ILE:HD12	1:E:393:HIS:NE2	2.19	0.57
1:E:280:ALA:HB2	1:E:296:VAL:HA	1.85	0.57
1:A:257:MSE:SE	1:B:257:MSE:HE2	2.54	0.57
1:A:266:GLY:HA3	1:B:295:GLU:HB3	1.87	0.57
1:D:153:ALA:HA	1:D:156:MSE:CE	2.34	0.57
1:D:162:LEU:HD21	1:D:282:LEU:HG	1.87	0.57
1:A:243:ASN:ND2	1:A:259:TRP:HE1	1.98	0.56
1:A:226:LEU:C	1:A:226:LEU:HD12	2.26	0.56
1:D:170:MSE:HE1	1:D:248:MSE:HE1	1.86	0.56
1:D:48:ILE:O	1:D:50:PRO:HD3	2.05	0.56
1:A:280:ALA:HB2	1:A:296:VAL:HA	1.88	0.56
1:E:172:PHE:CG	1:E:206:VAL:HB	2.41	0.55
1:E:115:LYS:NZ	1:E:205:ASP:O	2.40	0.55
1:A:128:ARG:HA	1:A:363:ILE:HG12	1.89	0.55
1:A:226:LEU:HB3	1:A:248:MSE:HG2	1.89	0.55
1:D:226:LEU:C	1:D:226:LEU:HD12	2.28	0.55
1:D:230:ARG:HD2	1:D:241:GLU:HB3	1.89	0.54
1:B:65:GLU:H	1:B:65:GLU:CD	2.11	0.54
1:D:156:MSE:HE1	1:D:292:LEU:HD21	1.89	0.54
1:B:271:GLN:HB3	1:B:286:ASP:CG	2.27	0.54
1:D:153:ALA:HA	1:D:156:MSE:HE3	1.90	0.54

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:D:16:MSE:HE1	1:D:29:HIS:CB	2.34	0.54
1:B:56:PHE:O	1:B:59:GLN:HG2	2.08	0.54
1:B:230:ARG:HG2	1:B:380:ALA:HB1	1.90	0.53
1:D:126:ASN:O	1:D:129:GLU:HB3	2.07	0.53
1:B:48:ILE:HG13	1:B:49:ASP:N	2.23	0.53
1:A:207:GLY:O	1:A:210:PRO:HD2	2.08	0.53
1:B:280:ALA:HB2	1:B:296:VAL:HA	1.91	0.53
1:E:226:LEU:C	1:E:226:LEU:HD12	2.29	0.53
1:A:98:TYR:CD1	1:A:123:GLN:HB3	2.43	0.53
1:D:179:ALA:C	3:D:549:HOH:O	2.47	0.53
1:E:395:HIS:HE1	3:E:655:HOH:O	1.90	0.53
1:D:351:ASP:OD1	1:D:354:ALA:HB3	2.08	0.53
1:D:209:HIS:HB3	1:D:271:GLN:HE22	1.75	0.52
1:B:224:LYS:HD3	1:B:251:GLU:HG3	1.91	0.52
1:E:170:MSE:HE1	1:E:248:MSE:HE1	1.92	0.52
1:D:280:ALA:HB2	1:D:296:VAL:HA	1.91	0.52
1:D:222:LYS:HE2	1:D:251:GLU:OE2	2.09	0.52
1:D:128:ARG:HA	1:D:363:ILE:HG12	1.91	0.52
1:B:92:THR:HB	1:B:93:PRO:CD	2.39	0.52
1:B:198:GLY:HA3	3:B:452:HOH:O	2.09	0.52
1:B:115:LYS:HB3	1:B:115:LYS:HZ2	1.75	0.51
1:E:128:ARG:HA	1:E:363:ILE:HG12	1.92	0.51
1:E:280:ALA:CB	1:E:296:VAL:HA	2.40	0.51
1:B:382:ASN:O	1:B:385:ILE:HG23	2.09	0.51
1:A:209:HIS:HB2	1:A:210:PRO:HD3	1.92	0.51
1:A:143:GLY:HA2	1:A:336:LEU:HD22	1.93	0.51
1:D:81:ARG:HB3	1:D:83:ASP:OD1	2.09	0.51
1:D:115:LYS:HB3	1:D:116:PRO:HD3	1.93	0.51
1:D:243:ASN:ND2	1:D:259:TRP:HE1	2.01	0.51
1:E:234:VAL:HB	1:E:237:ARG:HD2	1.92	0.51
1:A:339:ARG:HD3	1:A:359:ARG:O	2.11	0.51
1:D:282:LEU:HA	1:D:293:SER:O	2.11	0.50
1:B:170:MSE:HB3	1:B:210:PRO:HB2	1.91	0.50
1:E:243:ASN:HD22	1:E:261:SER:CB	2.24	0.50
1:A:209:HIS:HB3	1:A:271:GLN:HE22	1.76	0.50
1:B:198:GLY:N	3:B:452:HOH:O	2.45	0.50
1:D:295:GLU:HB3	1:E:266:GLY:HA3	1.94	0.50
1:A:263:VAL:HG23	1:B:167:MSE:HE1	1.93	0.50
1:B:72:LEU:O	1:B:76:GLU:HG2	2.12	0.50
1:B:226:LEU:C	1:B:226:LEU:HD12	2.33	0.49
1:B:110:HIS:HB3	1:B:344:MSE:HE2	1.94	0.49
1:A:227:MSE:SE	1:B:227:MSE:SE	3.31	0.49

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:154:ARG:NH2	1:A:359:ARG:HH11	2.11	0.49
1:B:44:GLY:HA2	1:B:74:MSE:SE	2.62	0.49
1:A:248:MSE:HE2	1:A:258:VAL:HG23	1.95	0.49
1:A:172:PHE:CG	1:A:206:VAL:HB	2.48	0.49
1:A:53:GLY:HA3	3:A:616:HOH:O	2.12	0.49
1:D:143:GLY:HA2	1:D:336:LEU:HD22	1.95	0.49
1:A:224:LYS:HD3	1:A:251:GLU:HG3	1.94	0.49
1:B:102:LYS:O	1:B:106:GLU:HG3	2.13	0.49
1:B:65:GLU:OE1	1:B:65:GLU:N	2.35	0.48
1:B:58:GLU:HA	1:B:58:GLU:OE1	2.13	0.48
1:D:373:VAL:O	1:D:377:VAL:HG23	2.13	0.48
1:D:227:MSE:SE	1:E:227:MSE:SE	3.31	0.48
1:E:98:TYR:CD1	1:E:123:GLN:HB3	2.49	0.48
1:E:157:ILE:HG12	1:E:162:LEU:HD22	1.96	0.48
1:A:210:PRO:HG3	1:A:271:GLN:HG3	1.97	0.47
1:A:56:PHE:CE2	1:A:60:LEU:HD11	2.49	0.47
1:E:16:MSE:HE1	1:E:30:ARG:HH11	1.80	0.47
1:A:340:PHE:O	1:A:344:MSE:HG3	2.14	0.47
1:A:280:ALA:CB	1:A:296:VAL:HA	2.45	0.47
1:B:200:SER:HB2	1:B:241:GLU:OE2	2.15	0.47
1:D:354:ALA:N	3:D:453:HOH:O	2.47	0.47
1:D:90:ILE:HD12	1:D:90:ILE:N	2.30	0.47
1:D:230:ARG:HG2	1:D:380:ALA:HB1	1.97	0.47
1:D:170:MSE:HB3	1:D:210:PRO:HB2	1.96	0.47
1:B:92:THR:HB	1:B:93:PRO:HD2	1.97	0.46
1:D:209:HIS:HB3	1:D:271:GLN:NE2	2.30	0.46
1:E:243:ASN:ND2	1:E:259:TRP:HE1	2.01	0.46
1:B:25:ILE:HD13	3:B:525:HOH:O	2.15	0.46
1:D:280:ALA:CB	1:D:296:VAL:HA	2.45	0.46
1:B:230:ARG:HD2	1:B:241:GLU:HB3	1.96	0.46
1:B:170:MSE:HE1	1:B:248:MSE:HE1	1.98	0.46
1:A:19:GLY:HA2	1:A:23:SER:CB	2.46	0.46
1:E:282:LEU:HA	1:E:293:SER:O	2.15	0.46
1:B:165:VAL:HG21	1:B:218:LEU:CD2	2.46	0.46
1:B:172:PHE:CG	1:B:206:VAL:HB	2.51	0.46
1:A:120:THR:OG1	1:A:123:GLN:HG3	2.16	0.46
1:D:317:ILE:HG13	1:D:318:ASP:N	2.31	0.45
1:D:115:LYS:HZ3	1:D:209:HIS:CD2	2.34	0.45
1:A:202:VAL:HG22	1:A:241:GLU:HG2	1.98	0.45
1:A:371:ARG:O	1:A:375:ARG:HG2	2.16	0.45
1:E:54:SER:O	1:E:58:GLU:HG3	2.16	0.45
1:D:271:GLN:HB3	1:D:286:ASP:CG	2.37	0.45

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:10:GLN:HE21	1:B:10:GLN:HB3	1.65	0.45
1:B:94:ASN:CG	1:B:116:PRO:HD2	2.37	0.45
1:D:89:SER:HA	1:D:112:VAL:O	2.16	0.45
1:A:271:GLN:HB3	1:A:286:ASP:CG	2.37	0.45
1:D:102:LYS:O	1:D:106:GLU:HG3	2.16	0.44
1:A:165:VAL:HG21	1:A:218:LEU:HD21	1.98	0.44
1:A:167:MSE:HE1	1:B:263:VAL:HG23	2.00	0.44
1:B:94:ASN:OD1	1:B:116:PRO:HD2	2.18	0.44
1:A:17:VAL:HG11	1:A:100:ILE:HD13	1.99	0.44
1:A:282:LEU:HA	1:A:293:SER:O	2.18	0.44
1:B:202:VAL:HG22	1:B:241:GLU:O	2.17	0.44
1:E:371:ARG:HD2	1:E:389:TYR:CZ	2.53	0.44
1:D:227:MSE:HE3	1:D:386:TRP:NE1	2.33	0.44
1:E:111:VAL:HG12	1:E:112:VAL:N	2.32	0.44
1:B:354:ALA:O	1:B:358:VAL:HG23	2.17	0.44
1:B:152:GLN:O	1:B:156:MSE:HG3	2.18	0.43
1:D:167:MSE:HE1	1:E:263:VAL:HG23	1.99	0.43
1:B:280:ALA:CB	1:B:296:VAL:HA	2.48	0.43
1:A:304:LEU:N	1:A:304:LEU:HD12	2.34	0.43
1:D:363:ILE:HG23	1:D:364:ASP:N	2.33	0.43
1:E:243:ASN:ND2	1:E:261:SER:HB2	2.33	0.43
1:D:234:VAL:HB	1:D:237:ARG:HD2	2.00	0.43
1:D:161:GLU:O	1:D:279:ARG:NH1	2.50	0.43
1:D:303:ILE:N	1:D:303:ILE:HD12	2.34	0.43
1:B:24:GLN:NE2	1:D:322:GLY:H	2.16	0.43
1:D:8:ILE:HG22	1:D:10:GLN:O	2.18	0.43
1:A:201:TYR:HB3	1:A:241:GLU:O	2.18	0.43
1:D:136:ARG:NH1	1:D:348:ASP:OD1	2.52	0.43
1:B:25:ILE:HD12	1:B:25:ILE:N	2.34	0.43
1:A:102:LYS:HG3	1:A:130:LEU:CD1	2.49	0.43
1:E:371:ARG:O	1:E:375:ARG:HG2	2.19	0.43
1:D:29:HIS:CE1	1:D:141:THR:HG21	2.54	0.43
1:D:120:THR:OG1	1:D:123:GLN:HG3	2.19	0.43
1:B:143:GLY:HA2	1:B:336:LEU:HD22	2.00	0.42
1:D:266:GLY:HA3	1:E:295:GLU:HB3	2.01	0.42
1:B:155:GLU:HG3	1:B:312:HIS:CG	2.54	0.42
1:D:115:LYS:HZ3	1:D:209:HIS:CG	2.38	0.42
1:B:230:ARG:HB2	1:B:241:GLU:CB	2.50	0.42
1:D:92:THR:HB	1:D:93:PRO:HD2	2.02	0.42
1:A:259:TRP:CE2	1:B:169:HIS:HE1	2.38	0.42
1:E:225:ARG:HE	1:E:249:GLU:CD	2.23	0.42
1:E:92:THR:HB	1:E:93:PRO:HD2	2.02	0.42

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:226:LEU:CB	1:A:248:MSE:HG2	2.50	0.42
1:B:81:ARG:HB3	1:B:83:ASP:OD1	2.20	0.42
1:B:234:VAL:HB	1:B:237:ARG:HD2	2.02	0.42
1:A:373:VAL:O	1:A:377:VAL:HG23	2.19	0.42
1:A:248:MSE:CE	1:A:258:VAL:HG23	2.50	0.41
1:A:209:HIS:HB3	1:A:271:GLN:NE2	2.34	0.41
1:B:96:THR:O	1:B:100:ILE:HG13	2.20	0.41
1:B:8:ILE:HD12	1:B:345:ASP:CG	2.40	0.41
1:E:157:ILE:HD12	1:E:217:MSE:HB3	2.02	0.41
1:D:8:ILE:HD12	1:D:341:ALA:HB1	2.02	0.41
1:B:224:LYS:HD2	1:B:224:LYS:HA	1.90	0.41
1:B:212:TYR:CE2	1:B:361:PRO:HG3	2.55	0.41
1:E:143:GLY:HA2	1:E:336:LEU:HD22	2.01	0.41
1:E:102:LYS:HG3	1:E:130:LEU:CD1	2.51	0.41
1:D:247:LEU:HD13	1:D:257:MSE:HE2	2.01	0.41
1:D:30:ARG:NH1	3:D:477:HOH:O	2.49	0.41
1:E:154:ARG:NH1	3:E:476:HOH:O	2.49	0.41
1:B:115:LYS:HD2	1:B:208:THR:OG1	2.19	0.41
1:E:89:SER:HA	1:E:112:VAL:O	2.19	0.41
1:E:385:ILE:CD1	1:E:393:HIS:NE2	2.84	0.41
1:B:149:LEU:HD23	1:B:149:LEU:HA	1.91	0.41
1:E:25:ILE:HD11	1:E:91:ALA:HB1	2.03	0.41
1:B:135:ASN:O	1:B:136:ARG:HD3	2.20	0.41
1:D:115:LYS:HB3	1:D:116:PRO:CD	2.50	0.41
1:D:171:GLN:O	1:D:271:GLN:HA	2.20	0.41
1:D:227:MSE:HE1	3:D:458:HOH:O	2.20	0.41
1:B:136:ARG:HA	1:B:136:ARG:HD3	1.83	0.41
1:E:304:LEU:HD12	1:E:304:LEU:N	2.35	0.41
1:D:96:THR:O	1:D:100:ILE:HG13	2.21	0.41
1:D:156:MSE:HE1	1:D:292:LEU:CD2	2.51	0.41
1:E:120:THR:OG1	1:E:123:GLN:HG3	2.21	0.41
1:E:121:VAL:O	1:E:125:GLU:HG2	2.20	0.41
1:D:156:MSE:CE	1:D:292:LEU:HD21	2.50	0.40
1:D:17:VAL:HG23	1:D:88:VAL:CG2	2.50	0.40
1:E:202:VAL:HG22	1:E:241:GLU:HG2	2.02	0.40
1:D:201:TYR:HB3	1:D:241:GLU:O	2.22	0.40
1:B:201:TYR:CD2	1:B:202:VAL:HG13	2.56	0.40
1:B:201:TYR:HB3	1:B:241:GLU:O	2.22	0.40
1:B:128:ARG:HA	1:B:363:ILE:HG12	2.02	0.40
1:E:382:ASN:O	1:E:385:ILE:HG23	2.22	0.40
1:A:268:MSE:HE3	1:A:268:MSE:HB3	1.96	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	370/398 (93%)	359 (97%)	10 (3%)	1 (0%)	50	39
1	B	360/398 (90%)	348 (97%)	11 (3%)	1 (0%)	50	39
1	D	354/398 (89%)	339 (96%)	12 (3%)	3 (1%)	27	14
1	E	370/398 (93%)	353 (95%)	16 (4%)	1 (0%)	50	39
All	All	1454/1592 (91%)	1399 (96%)	49 (3%)	6 (0%)	43	32

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	115	LYS
1	D	269	HIS
1	A	269	HIS
1	B	269	HIS
1	D	350	SER
1	E	269	HIS

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	297/303 (98%)	291 (98%)	6 (2%)	68	64
1	B	288/303 (95%)	283 (98%)	5 (2%)	73	71
1	D	286/303 (94%)	283 (99%)	3 (1%)	85	86
1	E	297/303 (98%)	291 (98%)	6 (2%)	68	64
All	All	1168/1212 (96%)	1148 (98%)	20 (2%)	73	71

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	GLN
1	A	72	LEU
1	A	115	LYS
1	A	172	PHE
1	A	269	HIS
1	A	397	HIS
1	B	10	GLN
1	B	65	GLU
1	B	72	LEU
1	B	172	PHE
1	B	269	HIS
1	D	172	PHE
1	D	269	HIS
1	D	351	ASP
1	E	24	GLN
1	E	115	LYS
1	E	172	PHE
1	E	269	HIS
1	E	353	GLN
1	E	397	HIS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (24) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	110	HIS
1	A	177	HIS
1	A	243	ASN
1	A	271	GLN
1	A	290	ASN
1	A	392	HIS
1	A	394	HIS
1	B	10	GLN
1	B	24	GLN
1	B	177	HIS
1	B	243	ASN
1	B	271	GLN
1	B	290	ASN
1	D	126	ASN
1	D	132	HIS
1	D	243	ASN
1	D	271	GLN
1	D	290	ASN
1	E	243	ASN

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Mol	Chain	Res	Type
1	E	271	GLN
1	E	290	ASN
1	E	392	HIS
1	E	394	HIS
1	E	395	HIS

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 4 ligands modelled in this entry, 4 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

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	374/398 (93%)	-0.27	9 (2%) 56 57	6, 13, 28, 36	0
1	B	364/398 (91%)	0.12	19 (5%) 26 25	7, 18, 34, 40	0
1	D	362/398 (90%)	0.03	9 (2%) 54 55	6, 17, 33, 40	0
1	E	374/398 (93%)	-0.29	6 (1%) 68 70	6, 12, 27, 37	0
All	All	1474/1592 (92%)	-0.10	43 (2%) 49 49	6, 15, 32, 40	0

All (43) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	51	ILE	5.6
1	B	22	GLN	4.9
1	B	8	ILE	4.8
1	B	48	ILE	4.7
1	D	8	ILE	4.3
1	D	48	ILE	3.8
1	B	9	PRO	3.8
1	B	352	THR	3.4
1	B	52	ARG	3.4
1	B	55	ALA	3.4
1	B	58	GLU	3.4
1	D	51	ILE	3.3
1	D	49	ASP	3.2
1	B	356	SER	3.2
1	D	9	PRO	3.2
1	B	50	PRO	3.2
1	B	353	GLN	3.1
1	D	115	LYS	3.0
1	A	48	ILE	3.0
1	A	235	ALA	2.9
1	B	350	SER	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	10	GLN	2.8
1	B	49	ASP	2.7
1	E	24	GLN	2.7
1	E	10	GLN	2.6
1	B	64	SER	2.6
1	A	51	ILE	2.6
1	B	10	GLN	2.6
1	B	129	GLU	2.5
1	E	235	ALA	2.4
1	E	22	GLN	2.4
1	D	356	SER	2.4
1	E	51	ILE	2.4
1	A	59	GLN	2.4
1	A	50	PRO	2.3
1	A	58	GLU	2.3
1	B	122	GLU	2.2
1	D	350	SER	2.1
1	A	22	GLN	2.1
1	E	58	GLU	2.1
1	D	64	SER	2.1
1	B	351	ASP	2.0
1	A	24	GLN	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(\AA^2)	Q<0.9
2	MG	A	399	1/1	0.13	0.86	29,29,29,29	0
2	MG	E	399	1/1	0.12	-0.48	30,30,30,30	0
2	MG	E	398	1/1	0.05	-3.16	17,17,17,17	0
2	MG	A	398	1/1	0.04	-3.90	19,19,19,19	0

6.5 Other polymers

There are no such residues in this entry.