



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 08:14 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 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

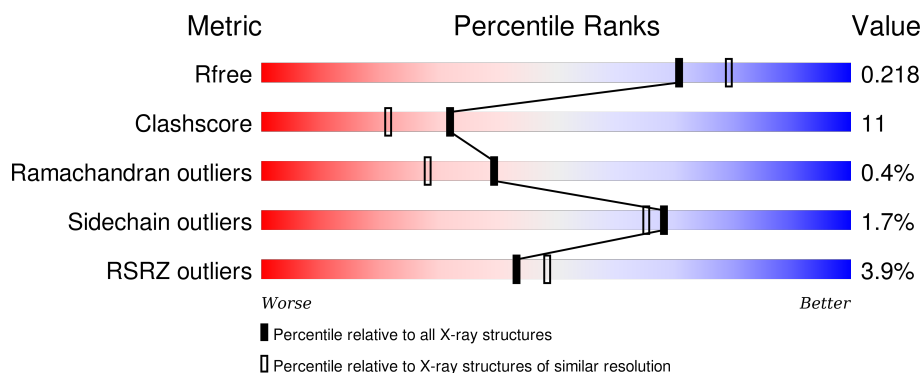
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.04 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}	91344	1192 (2.04-2.04)
Clashscore	102246	1269 (2.04-2.04)
Ramachandran outliers	100387	1258 (2.04-2.04)
Sidechain outliers	100360	1258 (2.04-2.04)
RSRZ outliers	91569	1194 (2.04-2.04)

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	398	<div> <div>3%</div> <div> <div></div> <div>75%</div> <div>18%</div> <div>• 6%</div> </div> </div>
1	B	398	<div> <div>6%</div> <div> <div></div> <div>69%</div> <div>21%</div> <div>• 9%</div> </div> </div>
1	D	398	<div> <div>4%</div> <div> <div></div> <div>69%</div> <div>22%</div> <div>• 9%</div> </div> </div>
1	E	398	<div> <div>2%</div> <div> <div></div> <div>75%</div> <div>19%</div> <div>• 6%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 12349 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 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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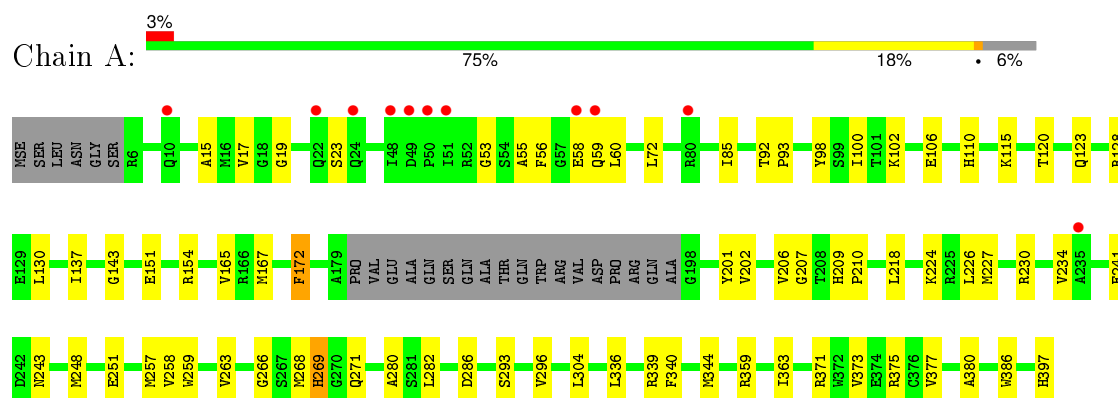
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	E	263	Total	O	0	0
			263	263		

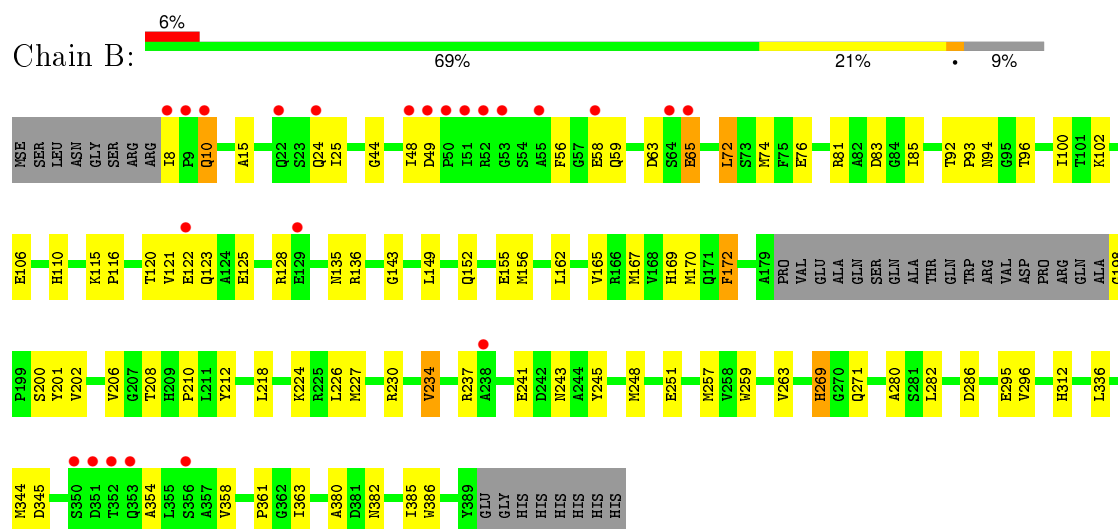
3 Residue-property plots [i](#)

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.

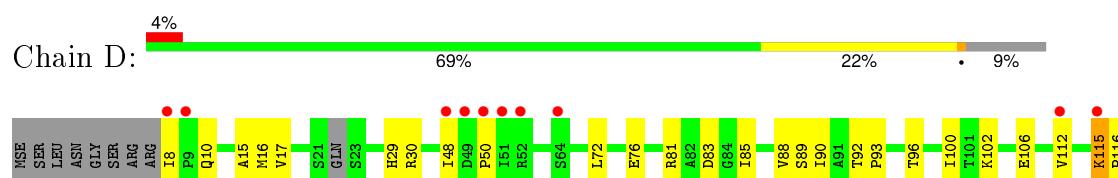
- Molecule 1: Oxidoreductase, Gfo/Idh/MocA family

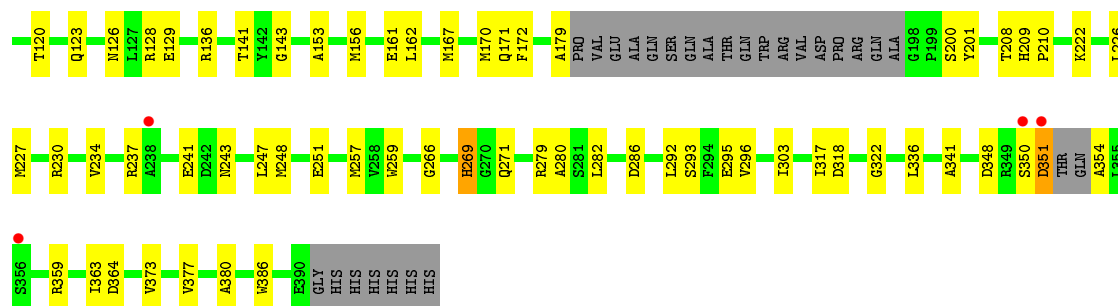


- Molecule 1: Oxidoreductase, Gfo/Idh/MocA family

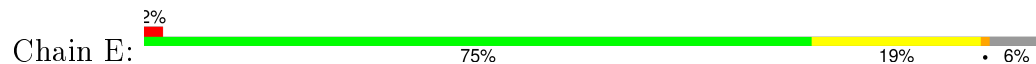


- Molecule 1: Oxidoreductase, Gfo/Idh/MocA family





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



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) 94.1 (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)	0.38 , 52.9	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	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.

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: 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 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	2914	0	2792	52	0
1	B	2818	0	2715	70	0
1	D	2802	0	2696	74	0
1	E	2914	0	2792	52	0
2	A	2	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
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

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 11.

All (237) 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: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	Interatomic distance (Å)	Clash overlap (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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:153:ALA:HA	1:D:156:MSE:HE3	1.90	0.54
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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
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

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
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 [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	370/398 (93%)	359 (97%)	10 (3%)	1 (0%)	46	36
1	B	360/398 (90%)	348 (97%)	11 (3%)	1 (0%)	46	36
1	D	354/398 (89%)	339 (96%)	12 (3%)	3 (1%)	24	12
1	E	370/398 (93%)	353 (95%)	16 (4%)	1 (0%)	46	36
All	All	1454/1592 (91%)	1399 (96%)	49 (3%)	6 (0%)	39	28

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 [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	297/303 (98%)	291 (98%)	6 (2%)	63	58
1	B	288/303 (95%)	283 (98%)	5 (2%)	68	65
1	D	286/303 (94%)	283 (99%)	3 (1%)	82	81
1	E	297/303 (98%)	291 (98%)	6 (2%)	63	58
All	All	1168/1212 (96%)	1148 (98%)	20 (2%)	68	65

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

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

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](#)

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.

No monomer is involved in short contacts.

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	361/398 (90%)	-0.21	11 (3%) 54 61	6, 12, 28, 36	0
1	B	351/398 (88%)	0.22	23 (6%) 22 24	7, 18, 34, 40	0
1	D	349/398 (87%)	0.12	14 (4%) 42 48	6, 16, 34, 40	0
1	E	361/398 (90%)	-0.24	7 (1%) 70 75	6, 12, 27, 37	0
All	All	1422/1592 (89%)	-0.03	55 (3%) 43 48	6, 14, 32, 40	0

All (55) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	51	ILE	6.7
1	B	48	ILE	5.7
1	B	22	GLN	5.0
1	B	8	ILE	4.7
1	B	9	PRO	4.6
1	D	8	ILE	4.4
1	D	48	ILE	4.3
1	D	51	ILE	4.0
1	B	52	ARG	3.9
1	B	58	GLU	3.9
1	B	352	THR	3.9
1	D	49	ASP	3.8
1	B	356	SER	3.7
1	B	353	GLN	3.7
1	B	50	PRO	3.6
1	A	48	ILE	3.6
1	B	55	ALA	3.6
1	D	9	PRO	3.5
1	B	49	ASP	3.4
1	E	24	GLN	3.4
1	A	235	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
1	D	115	LYS	3.2
1	E	51	ILE	3.0
1	B	350	SER	3.0
1	D	356	SER	3.0
1	E	235	ALA	2.9
1	E	10	GLN	2.8
1	A	51	ILE	2.8
1	B	64	SER	2.8
1	B	129	GLU	2.8
1	A	10	GLN	2.7
1	B	10	GLN	2.7
1	E	22	GLN	2.7
1	A	50	PRO	2.7
1	D	50	PRO	2.7
1	A	59	GLN	2.6
1	B	351	ASP	2.4
1	D	350	SER	2.4
1	A	58	GLU	2.4
1	B	53	GLY	2.4
1	D	238	ALA	2.3
1	E	58	GLU	2.3
1	A	24	GLN	2.3
1	B	122	GLU	2.3
1	D	351	ASP	2.2
1	A	22	GLN	2.2
1	B	24	GLN	2.1
1	D	64	SER	2.1
1	A	80	ARG	2.1
1	B	65	GLU	2.1
1	A	49	ASP	2.1
1	B	238	ALA	2.1
1	D	52	ARG	2.0
1	D	112	VAL	2.0
1	E	63	ASP	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 [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(\AA^2)	Q<0.9
2	MG	A	399	1/1	0.87	0.15	-	29,29,29,29	0
2	MG	A	398	1/1	0.97	0.05	-	19,19,19,19	0
2	MG	E	399	1/1	0.92	0.14	-	30,30,30,30	0
2	MG	E	398	1/1	0.94	0.05	-	17,17,17,17	0

6.5 Other polymers [i](#)

There are no such residues in this entry.