



Full wwPDB X-ray Structure Validation Report ⓘ

Jun 17, 2024 – 10:27 PM EDT

PDB ID : 5YXY
Title : Crystal structure of the HyhL-HypA complex (form I)
Authors : Kwon, S.; Watanabe, S.; Nishitani, Y.; Miki, K.
Deposited on : 2017-12-07
Resolution : 3.30 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

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

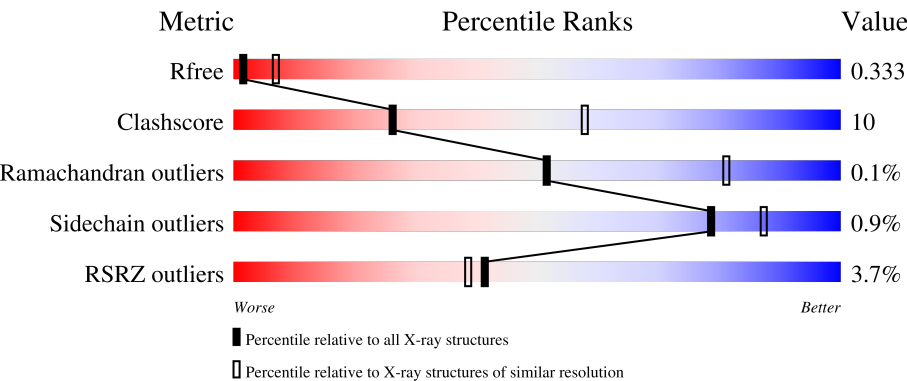
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 3.30 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}	130704	1149 (3.34-3.26)
Clashscore	141614	1205 (3.34-3.26)
Ramachandran outliers	138981	1183 (3.34-3.26)
Sidechain outliers	138945	1182 (3.34-3.26)
RSRZ outliers	127900	1115 (3.34-3.26)

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 of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. 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	428	<div><div>%</div><div>76%18%6%</div></div>
1	B	428	<div><div>6%</div><div>76%16%8%</div></div>
1	C	428	<div><div>%</div><div>74%20%• 5%</div></div>
2	D	139	<div><div>4%</div><div>77%12%• 11%</div></div>
2	E	139	<div><div>7%</div><div>75%13%• 12%</div></div>

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Mol	Chain	Length	Quality of chain
2	F	139	<div><div></div><div>6%</div><div>78%</div><div>9%</div><div>•</div><div>13%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 11198 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 Cytosolic NiFe-hydrogenase, alpha subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	404	Total	C	N	O	S	0	0	0
			2851	1814	476	545	16			
1	B	394	Total	C	N	O	S	0	0	0
			2697	1718	454	510	15			
1	C	406	Total	C	N	O	S	0	0	0
			2940	1887	492	546	15			

- Molecule 2 is a protein called Probable hydrogenase nickel incorporation protein HypA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	D	124	Total	C	N	O	S	0	0	0
			917	587	155	170	5			
2	E	123	Total	C	N	O	S	0	0	0
			925	594	153	173	5			
2	F	121	Total	C	N	O	S	0	0	0
			865	555	139	166	5			

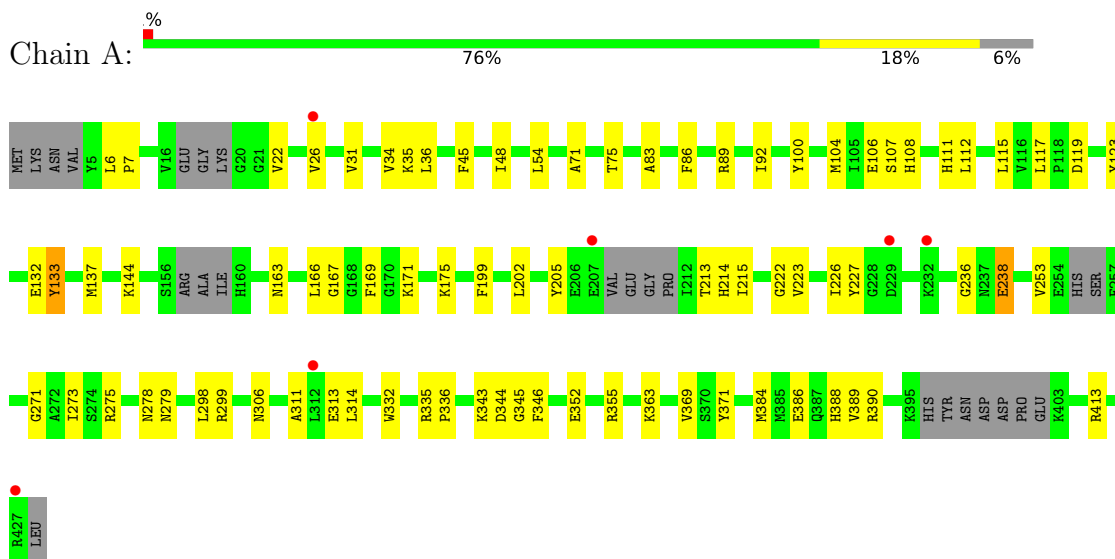
- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	1	Total	Zn	0	0
			1	1		
3	E	1	Total	Zn	0	0
			1	1		
3	F	1	Total	Zn	0	0
			1	1		

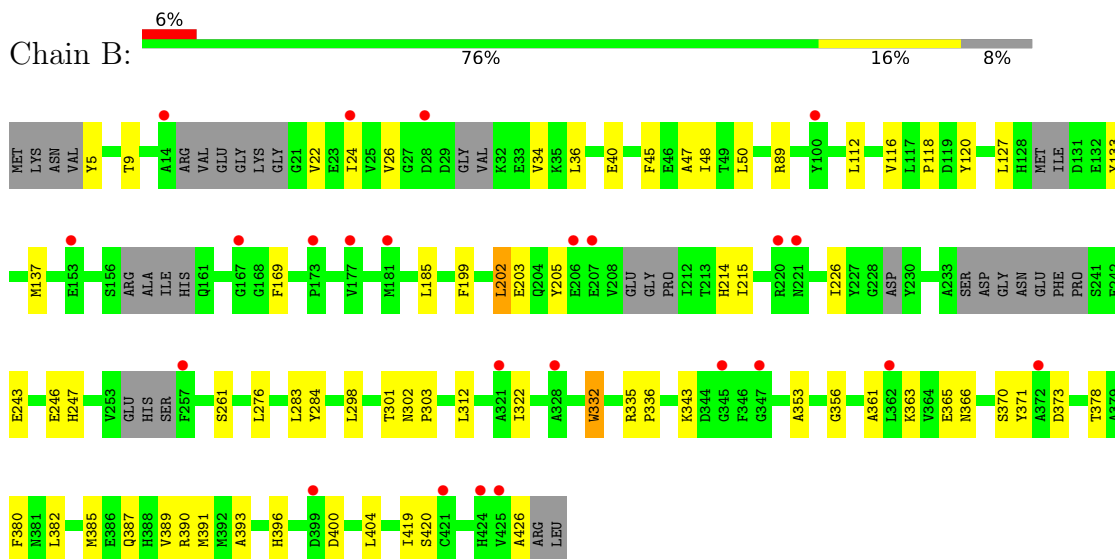
3 Residue-property plots [i](#)

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

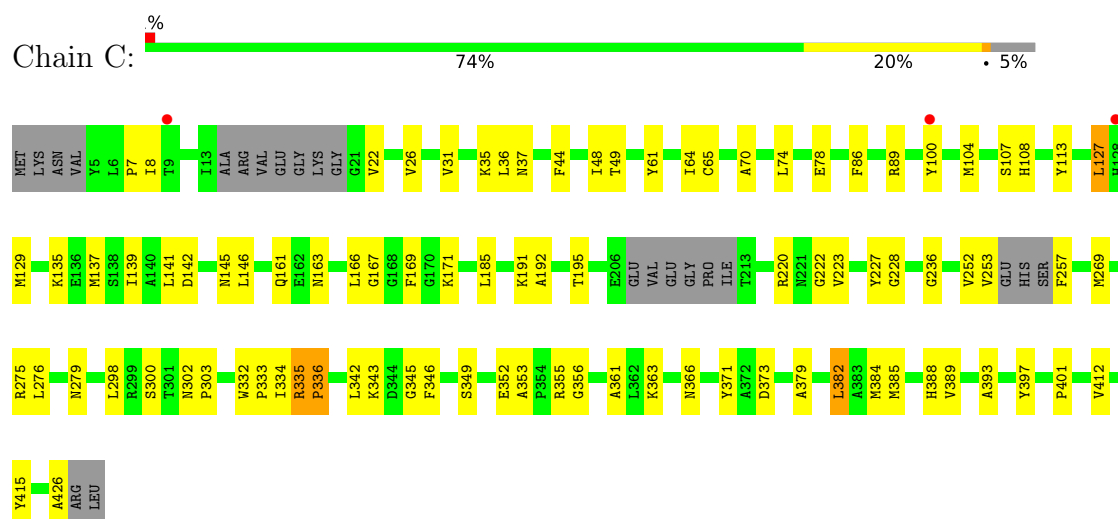
- Molecule 1: Cytosolic NiFe-hydrogenase, alpha subunit



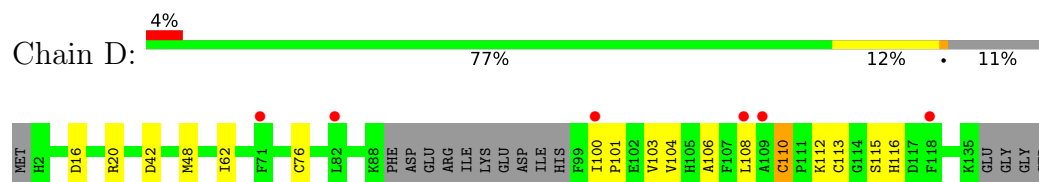
- Molecule 1: Cytosolic NiFe-hydrogenase, alpha subunit



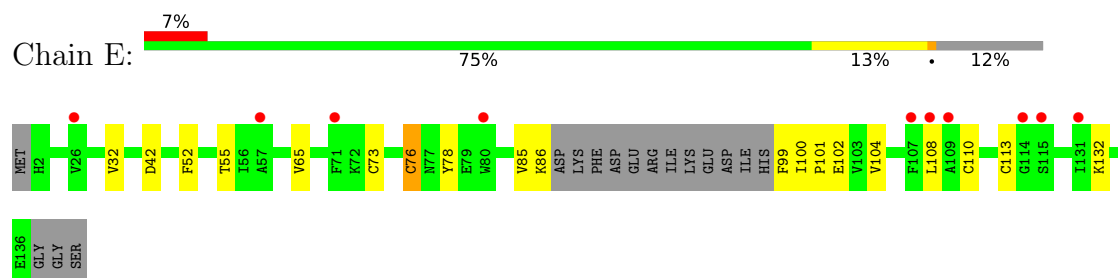
- Molecule 1: Cytosolic NiFe-hydrogenase, alpha subunit



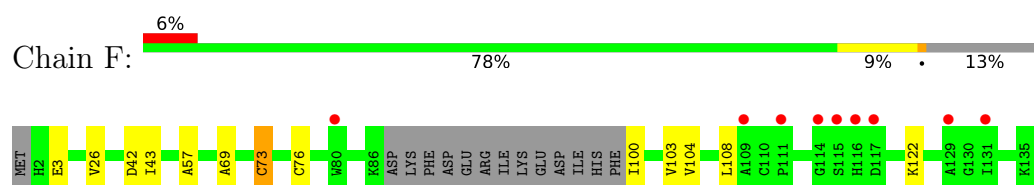
- Molecule 2: Probable hydrogenase nickel incorporation protein HypA



- Molecule 2: Probable hydrogenase nickel incorporation protein HypA



- Molecule 2: Probable hydrogenase nickel incorporation protein HypA



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 2	Depositor
Cell constants a, b, c, α , β , γ	245.31 Å 261.43 Å 134.87 Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.22 – 3.30 48.22 – 3.30	Depositor EDS
% Data completeness (in resolution range)	99.0 (48.22-3.30) 99.0 (48.22-3.30)	Depositor EDS
R_{merge}	0.15	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.97 (at 3.33 Å)	Xtriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.305 , 0.321 0.318 , 0.333	Depositor DCC
R_{free} test set	3183 reflections (4.91%)	wwPDB-VP
Wilson B-factor (Å ²)	87.3	Xtriage
Anisotropy	0.701	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.27 , 34.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.85	EDS
Total number of atoms	11198	wwPDB-VP
Average B, all atoms (Å ²)	50.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.90% 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, 0.333 respectively for untwinned datasets, and 0.375, 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: ZN

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	0.27	0/2898	0.48	0/3953
1	B	0.27	0/2741	0.46	1/3743 (0.0%)
1	C	0.33	2/2996 (0.1%)	0.51	2/4079 (0.0%)
2	D	0.27	0/933	0.46	0/1267
2	E	0.24	0/941	0.42	0/1276
2	F	0.24	0/881	0.42	0/1203
All	All	0.28	2/11390 (0.0%)	0.47	3/15521 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	401	PRO	N-CD	-7.19	1.37	1.47
1	C	336	PRO	N-CD	5.18	1.55	1.47

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	382	LEU	CA-CB-CG	5.45	127.84	115.30
1	C	335	ARG	C-N-CD	5.45	139.84	128.40
1	B	332	TRP	C-N-CD	5.43	139.80	128.40

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	2851	0	2573	58	0
1	B	2697	0	2341	54	0
1	C	2940	0	2746	63	0
2	D	917	0	850	19	0
2	E	925	0	870	24	0
2	F	865	0	760	9	0
3	D	1	0	0	0	0
3	E	1	0	0	0	0
3	F	1	0	0	0	0
All	All	11198	0	10140	222	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 10.

All (222) 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)
2:E:76:CYS:HB3	2:E:113:CYS:SG	1.57	1.43
2:E:76:CYS:CB	2:E:113:CYS:SG	2.27	1.21
1:A:123:TYR:CE2	1:A:132:GLU:OE1	1.92	1.21
1:B:26:VAL:O	1:B:426:ALA:CB	2.03	1.07
1:A:238:GLU:OE1	1:A:238:GLU:N	1.88	1.06
2:E:76:CYS:SG	2:E:78:TYR:HB3	1.95	1.05
2:E:76:CYS:SG	2:E:78:TYR:CB	2.44	1.04
1:A:123:TYR:HE2	1:A:132:GLU:OE1	1.40	0.98
1:B:26:VAL:O	1:B:426:ALA:HB1	1.64	0.98
1:A:384:MET:HE3	1:A:388:HIS:NE2	1.84	0.92
1:C:332:TRP:CD2	1:C:333:PRO:HA	2.05	0.91
1:A:236:GLY:O	1:A:278:ASN:ND2	2.04	0.91
2:E:76:CYS:SG	2:E:78:TYR:HB2	2.12	0.90
1:A:346:PHE:HD1	1:A:363:LYS:HB2	1.38	0.89
2:E:85:VAL:O	2:E:86:LYS:HG2	1.74	0.88
1:A:384:MET:CE	1:A:388:HIS:NE2	2.36	0.88
2:D:104:VAL:CG1	2:D:108:LEU:HB2	2.05	0.86
2:D:76:CYS:HB3	2:D:113:CYS:SG	2.21	0.81
1:C:269:MET:HB2	1:C:384:MET:HG2	1.64	0.79
2:E:99:PHE:C	2:E:101:PRO:HD3	2.07	0.75
1:A:106:GLU:HG3	1:A:144:LYS:HD2	1.69	0.75
1:B:363:LYS:HB3	1:B:371:TYR:HB3	1.69	0.75
1:B:203:GLU:HG2	1:B:205:TYR:H	1.52	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:169:PHE:H	1:A:336:PRO:HB3	1.51	0.74
1:A:205:TYR:O	1:A:213:THR:HA	1.87	0.74
1:B:246:GLU:HG2	1:B:247:HIS:H	1.52	0.74
1:C:363:LYS:HB3	1:C:371:TYR:HB3	1.70	0.73
2:E:76:CYS:HB2	2:E:113:CYS:SG	2.29	0.72
1:B:169:PHE:H	1:B:336:PRO:HB3	1.53	0.72
1:A:363:LYS:HB3	1:A:371:TYR:HB3	1.71	0.71
1:A:346:PHE:CD1	1:A:363:LYS:HB2	2.25	0.71
2:D:104:VAL:HG12	2:D:108:LEU:HB2	1.72	0.71
2:E:104:VAL:HG13	2:E:108:LEU:HB2	1.73	0.70
1:B:5:TYR:HA	2:E:132:LYS:HD3	1.74	0.69
1:A:107:SER:OG	1:A:355:ARG:NH1	2.25	0.69
1:C:393:ALA:O	1:C:397:TYR:N	2.26	0.69
2:E:100:ILE:N	2:E:101:PRO:HD3	2.08	0.68
1:C:275:ARG:HD3	1:C:352:GLU:HG2	1.74	0.68
2:D:115:SER:O	2:D:116:HIS:HB2	1.91	0.68
1:A:26:VAL:HG12	1:A:31:VAL:HG12	1.76	0.67
1:C:332:TRP:CG	1:C:333:PRO:HA	2.28	0.67
1:B:246:GLU:HG2	1:B:247:HIS:N	2.10	0.67
2:D:104:VAL:CG1	2:D:108:LEU:HD22	2.25	0.67
1:A:215:ILE:HD12	1:A:299:ARG:HE	1.59	0.67
1:C:141:LEU:O	1:C:145:ASN:ND2	2.27	0.67
1:C:142:ASP:OD2	1:C:191:LYS:NZ	2.27	0.67
1:A:123:TYR:CZ	1:A:132:GLU:OE1	2.48	0.67
1:A:343:LYS:HG2	1:A:344:ASP:N	2.10	0.66
1:C:100:TYR:OH	1:C:352:GLU:O	2.05	0.65
2:D:104:VAL:HG13	2:D:108:LEU:HD22	1.78	0.65
2:E:85:VAL:O	2:E:86:LYS:CG	2.44	0.65
1:A:384:MET:HE2	1:A:388:HIS:NE2	2.12	0.65
1:C:127:LEU:HD23	1:C:129:MET:H	1.61	0.65
2:E:100:ILE:N	2:E:101:PRO:CD	2.60	0.65
1:B:215:ILE:HD11	1:B:301:THR:OG1	1.97	0.64
1:C:333:PRO:O	1:C:334:ILE:HG22	1.97	0.64
2:D:104:VAL:HG13	2:D:108:LEU:HB2	1.80	0.64
1:C:220:ARG:NH1	1:C:300:SER:O	2.30	0.64
2:E:76:CYS:HG	2:E:78:TYR:HB3	1.64	0.63
1:C:223:VAL:O	1:C:279:ASN:ND2	2.31	0.63
1:B:276:LEU:HD21	1:B:298:LEU:HD13	1.80	0.62
1:B:363:LYS:HG2	1:B:370:SER:HB3	1.81	0.62
1:A:100:TYR:OH	1:A:352:GLU:O	2.07	0.62
1:A:89:ARG:NH2	1:A:335:ARG:O	2.33	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:132:GLU:HG2	1:A:133:TYR:HB2	1.82	0.61
1:C:334:ILE:O	1:C:334:ILE:HG23	2.00	0.61
2:D:103:VAL:HG12	2:D:103:VAL:O	2.01	0.60
1:B:22:VAL:HG13	1:B:36:LEU:HB3	1.82	0.60
1:C:86:PHE:CE2	1:C:167:GLY:HA2	2.37	0.60
2:F:73:CYS:SG	2:F:76:CYS:HB2	2.41	0.59
1:B:89:ARG:NH2	1:B:335:ARG:O	2.34	0.59
1:A:275:ARG:NH1	1:A:313:GLU:OE1	2.37	0.58
1:A:386:GLU:O	1:A:389:VAL:HG22	2.03	0.58
1:C:26:VAL:O	1:C:426:ALA:HA	2.04	0.58
1:C:345:GLY:O	1:C:363:LYS:HA	2.04	0.58
2:E:76:CYS:CA	2:E:113:CYS:SG	2.92	0.57
1:B:133:TYR:CE2	1:B:137:MET:HG2	2.40	0.57
1:C:107:SER:OG	1:C:355:ARG:NH1	2.37	0.57
1:A:89:ARG:H	1:A:92:ILE:HD11	1.70	0.57
1:C:161:GLN:OE1	1:C:161:GLN:N	2.38	0.57
1:A:273:ILE:HG12	1:A:306:ASN:HA	1.86	0.56
1:B:45:PHE:HA	1:B:48:ILE:HG12	1.87	0.56
1:C:192:ALA:O	1:C:195:THR:OG1	2.16	0.56
1:C:127:LEU:HD13	1:C:137:MET:HE2	1.88	0.56
1:A:45:PHE:HA	1:A:48:ILE:HG12	1.88	0.56
1:C:104:MET:SD	1:C:108:HIS:NE2	2.79	0.55
1:B:419:ILE:HG13	1:B:420:SER:N	2.21	0.55
1:C:361:ALA:HB3	1:C:373:ASP:HB3	1.87	0.55
1:B:390:ARG:HA	1:B:393:ALA:HB3	1.88	0.55
2:E:104:VAL:HG22	2:E:108:LEU:HD22	1.88	0.55
1:B:343:LYS:HG2	1:B:366:ASN:H	1.71	0.55
2:E:85:VAL:C	2:E:86:LYS:HG2	2.26	0.55
1:A:54:LEU:HD22	1:A:166:LEU:HD22	1.88	0.54
1:C:44:PHE:CZ	1:C:48:ILE:HD11	2.42	0.54
1:C:222:GLY:O	1:C:236:GLY:N	2.33	0.54
1:C:275:ARG:HH21	1:C:352:GLU:HG3	1.71	0.54
1:B:26:VAL:O	1:B:426:ALA:HB3	2.01	0.54
1:A:223:VAL:O	1:A:279:ASN:ND2	2.40	0.54
1:C:146:LEU:HD21	1:C:185:LEU:HD12	1.91	0.53
1:A:83:ALA:O	1:A:345:GLY:HA3	2.09	0.53
1:A:133:TYR:HD1	1:A:137:MET:HG3	1.73	0.53
1:A:384:MET:HE3	1:A:388:HIS:CD2	2.42	0.53
1:B:215:ILE:HG13	1:B:301:THR:HB	1.91	0.53
1:B:185:LEU:HD22	1:B:322:ILE:HB	1.90	0.52
1:C:343:LYS:HG3	1:C:366:ASN:H	1.72	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:TYR:CD1	1:A:137:MET:HG3	2.45	0.52
2:E:101:PRO:O	2:E:102:GLU:HB2	2.08	0.52
1:C:22:VAL:HG13	1:C:36:LEU:HB3	1.91	0.52
1:C:385:MET:HE1	1:C:412:VAL:HG13	1.91	0.52
1:C:169:PHE:H	1:C:336:PRO:HB3	1.75	0.51
1:C:37:ASN:OD1	1:C:37:ASN:N	2.43	0.51
2:F:3:GLU:HG2	2:F:43:ILE:HB	1.90	0.51
1:B:396:HIS:O	1:B:396:HIS:ND1	2.42	0.51
1:C:61:TYR:O	1:C:64:ILE:HD12	2.10	0.51
1:C:379:ALA:O	1:C:382:LEU:HG	2.10	0.51
1:A:214:HIS:O	1:A:390:ARG:NH2	2.43	0.51
2:D:16:ASP:OD2	2:D:20:ARG:NH2	2.44	0.50
1:B:215:ILE:HG13	1:B:301:THR:O	2.11	0.50
1:C:139:ILE:HD12	1:C:191:LYS:HB3	1.94	0.50
1:C:276:LEU:HD13	1:C:298:LEU:HD13	1.94	0.50
1:A:35:LYS:HD3	1:A:253:VAL:HA	1.94	0.50
1:C:353:ALA:N	1:C:356:GLY:O	2.35	0.49
1:B:215:ILE:HG13	1:B:301:THR:CB	2.41	0.49
2:D:100:ILE:N	2:D:101:PRO:CD	2.76	0.49
1:B:343:LYS:HE3	1:B:365:GLU:HG3	1.94	0.49
1:B:353:ALA:N	1:B:356:GLY:O	2.39	0.49
2:E:76:CYS:N	2:E:113:CYS:SG	2.86	0.49
1:A:104:MET:SD	1:A:108:HIS:NE2	2.86	0.48
1:C:70:ALA:HB2	1:C:161:GLN:NE2	2.27	0.48
1:C:191:LYS:O	1:C:195:THR:HG23	2.13	0.48
1:B:261:SER:OG	1:B:380:PHE:HA	2.14	0.48
1:B:120:TYR:CZ	1:B:202:LEU:HD21	2.49	0.48
1:C:44:PHE:CE2	2:F:43:ILE:HG12	2.48	0.48
2:D:104:VAL:O	2:D:106:ALA:N	2.47	0.48
1:A:311:ALA:O	1:A:314:LEU:HB2	2.14	0.48
1:B:215:ILE:CD1	1:B:301:THR:OG1	2.61	0.47
1:C:343:LYS:O	1:C:345:GLY:N	2.47	0.47
1:B:116:VAL:HG21	1:B:199:PHE:CZ	2.49	0.47
1:A:71:ALA:O	1:A:75:THR:HG22	2.14	0.47
2:D:76:CYS:CB	2:D:113:CYS:SG	2.95	0.47
2:D:104:VAL:HG13	2:D:108:LEU:HD13	1.96	0.47
2:F:26:VAL:HG12	2:F:57:ALA:HA	1.95	0.47
2:D:104:VAL:HG11	2:D:108:LEU:HD22	1.95	0.47
1:A:271:GLY:O	1:A:275:ARG:HD3	2.16	0.46
1:C:113:TYR:OH	1:C:195:THR:HG22	2.14	0.46
1:A:215:ILE:HD12	1:A:299:ARG:NE	2.29	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:47:ALA:HB1	1:C:7:PRO:HG2	1.96	0.46
1:B:24:ILE:HG12	1:B:34:VAL:HG23	1.97	0.46
2:D:42:ASP:OD1	2:D:42:ASP:N	2.49	0.46
2:E:52:PHE:O	2:E:55:THR:HG22	2.16	0.46
1:B:50:LEU:HD23	1:C:8:ILE:HA	1.98	0.45
1:B:385:MET:O	1:B:389:VAL:HG23	2.17	0.45
1:C:385:MET:O	1:C:389:VAL:HG23	2.16	0.45
1:A:369:VAL:O	1:B:9:THR:HG21	2.16	0.45
1:B:214:HIS:O	1:B:215:ILE:C	2.55	0.45
1:C:49:THR:HG21	1:C:61:TYR:CE1	2.52	0.45
1:A:89:ARG:HH22	1:A:335:ARG:H	1.64	0.45
1:C:166:LEU:HD12	1:C:166:LEU:H	1.82	0.45
2:E:73:CYS:HB2	2:E:110:CYS:HB2	1.99	0.45
1:B:120:TYR:CE2	1:B:202:LEU:HD21	2.52	0.44
1:B:133:TYR:CD2	1:B:133:TYR:O	2.70	0.44
2:D:104:VAL:CG1	2:D:108:LEU:CB	2.89	0.44
2:D:110:CYS:SG	2:D:112:LYS:N	2.85	0.44
1:C:135:LYS:O	1:C:139:ILE:HG12	2.17	0.44
1:B:283:LEU:HD11	1:B:312:LEU:HD21	1.99	0.44
1:B:283:LEU:HD23	1:B:284:TYR:N	2.32	0.44
1:C:89:ARG:NH2	1:C:335:ARG:O	2.48	0.44
1:C:253:VAL:HB	1:C:257:PHE:CE1	2.53	0.44
1:B:302:ASN:HA	1:B:303:PRO:HD3	1.79	0.44
1:A:163:ASN:ND2	1:A:171:LYS:O	2.29	0.44
1:C:388:HIS:ND1	1:C:415:TYR:OH	2.32	0.43
1:B:133:TYR:HE2	1:B:137:MET:HG2	1.81	0.43
2:F:100:ILE:HG22	2:F:100:ILE:O	2.18	0.43
1:A:119:ASP:OD2	1:A:413:ARG:NH1	2.51	0.43
1:A:226:ILE:CG1	1:A:227:TYR:N	2.81	0.43
1:B:243:GLU:HG2	1:B:243:GLU:O	2.18	0.43
1:A:386:GLU:O	1:A:389:VAL:CG2	2.66	0.43
2:E:100:ILE:HG22	2:E:100:ILE:O	2.17	0.43
1:C:26:VAL:HA	1:C:31:VAL:HA	1.99	0.43
1:C:31:VAL:HG11	1:C:393:ALA:HB1	2.00	0.43
1:C:78:GLU:OE1	1:C:349:SER:OG	2.36	0.43
2:E:42:ASP:N	2:E:42:ASP:OD1	2.51	0.43
2:F:42:ASP:OD1	2:F:42:ASP:N	2.50	0.43
1:A:175:LYS:NZ	1:A:332:TRP:O	2.41	0.43
1:B:343:LYS:HE2	1:B:343:LYS:HB3	1.83	0.43
1:C:74:LEU:HD13	1:C:100:TYR:HB2	2.00	0.42
2:F:100:ILE:HA	2:F:103:VAL:HB	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:ARG:HA	1:A:336:PRO:HD3	1.80	0.42
1:C:303:PRO:HG3	1:C:388:HIS:CE1	2.54	0.42
1:A:273:ILE:HD13	1:A:273:ILE:HA	1.87	0.42
1:B:378:THR:O	1:B:382:LEU:HG	2.19	0.42
2:D:104:VAL:HG13	2:D:108:LEU:CD2	2.49	0.42
1:C:382:LEU:O	1:C:385:MET:HG2	2.20	0.42
1:A:86:PHE:CE2	1:A:167:GLY:HA2	2.55	0.42
1:A:34:VAL:HG21	1:A:389:VAL:HG21	2.02	0.42
1:B:243:GLU:OE1	1:B:243:GLU:N	2.53	0.42
1:B:361:ALA:HB3	1:B:373:ASP:HB3	2.02	0.42
1:C:346:PHE:HD1	1:C:363:LYS:HB2	1.84	0.42
1:B:112:LEU:HD13	1:B:199:PHE:CE1	2.55	0.42
2:D:48:MET:HE2	2:D:62:ILE:HG12	2.01	0.41
1:A:199:PHE:CD2	1:A:202:LEU:HD21	2.55	0.41
1:B:118:PRO:HG3	1:B:127:LEU:HD11	2.02	0.41
1:C:227:TYR:O	1:C:228:GLY:C	2.57	0.41
2:E:32:VAL:HG22	2:E:65:VAL:HB	2.01	0.41
1:B:203:GLU:HG2	1:B:205:TYR:N	2.28	0.41
1:C:302:ASN:HA	1:C:303:PRO:HD3	1.82	0.41
1:A:117:LEU:HD23	1:A:117:LEU:HA	1.88	0.41
1:B:112:LEU:O	1:B:116:VAL:HG22	2.20	0.41
1:B:387:GLN:O	1:B:391:MET:HG3	2.20	0.41
1:B:400:ASP:HB2	1:B:404:LEU:CD2	2.51	0.41
2:F:104:VAL:O	2:F:108:LEU:HB2	2.21	0.41
1:A:22:VAL:HG12	1:A:36:LEU:HB2	2.03	0.41
1:C:35:LYS:NZ	1:C:252:VAL:O	2.54	0.41
1:A:112:LEU:HB3	1:A:199:PHE:CZ	2.56	0.41
1:C:163:ASN:OD1	1:C:171:LYS:N	2.52	0.41
1:A:6:LEU:HA	1:A:7:PRO:HD3	1.80	0.40
1:A:111:HIS:O	1:A:115:LEU:HB2	2.22	0.40
1:A:222:GLY:O	1:A:236:GLY:N	2.54	0.40
2:F:69:ALA:HA	2:F:122:LYS:O	2.21	0.40
1:A:343:LYS:CG	1:A:344:ASP:N	2.80	0.40
1:B:40:GLU:N	1:B:40:GLU:OE1	2.55	0.40
1:C:342:LEU:HD23	1:C:342:LEU:HA	1.98	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	392/428 (92%)	384 (98%)	8 (2%)	0	100	100
1	B	376/428 (88%)	366 (97%)	8 (2%)	2 (0%)	29	61
1	C	398/428 (93%)	388 (98%)	10 (2%)	0	100	100
2	D	120/139 (86%)	115 (96%)	5 (4%)	0	100	100
2	E	119/139 (86%)	118 (99%)	1 (1%)	0	100	100
2	F	117/139 (84%)	116 (99%)	1 (1%)	0	100	100
All	All	1522/1701 (90%)	1487 (98%)	33 (2%)	2 (0%)	51	81

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	332	TRP
1	B	226	ILE

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	255/361 (71%)	252 (99%)	3 (1%)	71	83
1	B	222/361 (62%)	221 (100%)	1 (0%)	88	93
1	C	273/361 (76%)	271 (99%)	2 (1%)	84	90
2	D	85/114 (75%)	84 (99%)	1 (1%)	71	83
2	E	88/114 (77%)	87 (99%)	1 (1%)	73	85

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	F	76/114 (67%)	75 (99%)	1 (1%)	69	82
All	All	999/1425 (70%)	990 (99%)	9 (1%)	78	87

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	133	TYR
1	A	238	GLU
1	A	298	LEU
1	B	202	LEU
1	C	65	CYS
1	C	127	LEU
2	D	110	CYS
2	E	76	CYS
2	F	73	CYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	398	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no monosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 3 ligands modelled in this entry, 3 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 [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	404/428 (94%)	0.05	6 (1%) 73 72	19, 39, 63, 86	0
1	B	394/428 (92%)	0.28	24 (6%) 21 20	31, 69, 93, 141	0
1	C	406/428 (94%)	0.14	3 (0%) 87 88	20, 40, 73, 90	0
2	D	124/139 (89%)	0.32	6 (4%) 30 28	20, 41, 62, 72	0
2	E	123/139 (88%)	0.39	10 (8%) 12 11	27, 55, 85, 102	0
2	F	121/139 (87%)	0.34	9 (7%) 14 14	35, 54, 70, 84	0
All	All	1572/1701 (92%)	0.20	58 (3%) 41 38	19, 48, 84, 141	0

All (58) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	229	ASP	7.4
2	D	82	LEU	4.5
1	B	206	GLU	4.4
2	F	111	PRO	4.2
1	A	207	GLU	3.8
1	B	153	GLU	3.7
2	E	71	PHE	3.6
2	D	109	ALA	3.4
1	B	328	ALA	3.3
2	D	108	LEU	3.3
1	B	372	ALA	3.3
2	D	118	PHE	3.3
1	B	221	ASN	3.1
1	B	220	ARG	3.1
1	B	424	HIS	3.1
1	B	362	LEU	3.0
2	E	114	GLY	2.9
1	B	24	ILE	2.8
1	B	28	ASP	2.8

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Mol	Chain	Res	Type	RSRZ
1	B	181	MET	2.8
1	B	425	VAL	2.7
2	F	109	ALA	2.7
2	E	108	LEU	2.6
1	B	345	GLY	2.6
2	D	100	ILE	2.6
2	F	114	GLY	2.6
2	D	71	PHE	2.6
1	C	100	TYR	2.5
2	F	129	ALA	2.5
2	F	80	TRP	2.5
1	B	347	GLY	2.5
1	B	100	TYR	2.5
1	B	14	ALA	2.4
1	B	173	PRO	2.4
2	E	26	VAL	2.4
1	C	128	HIS	2.4
1	B	257	PHE	2.3
2	E	107	PHE	2.3
2	F	116	HIS	2.3
1	B	167	GLY	2.3
2	F	117	ASP	2.3
1	B	399	ASP	2.2
1	A	427	ARG	2.2
2	F	115	SER	2.2
2	E	80	TRP	2.2
1	A	232	LYS	2.2
1	A	26	VAL	2.2
2	E	57	ALA	2.1
1	B	177	VAL	2.1
2	E	131	ILE	2.1
1	B	321	ALA	2.1
1	A	312	LEU	2.0
2	F	131	ILE	2.0
1	B	421	CYS	2.0
1	B	207	GLU	2.0
2	E	109	ALA	2.0
2	E	115	SER	2.0
1	C	9	THR	2.0

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides 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. 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	B-factors(\AA^2)	Q<0.9
3	ZN	E	701	1/1	0.86	0.05	75,75,75,75	0
3	ZN	D	701	1/1	0.93	0.12	43,43,43,43	0
3	ZN	F	701	1/1	0.98	0.06	43,43,43,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.