Full wwPDB X-ray Structure Validation Report ⓘ

Nov 19, 2023 – 08:27 PM JST

PDB ID	:	7BQW
Title	:	Crystal structure of Methionine gamma-lyase from Fusobacterium nucleatum
Authors	:	Lan, J.; Chen, Y.; Liu, W.; Xu, Y.
Deposited on	:	2020-03-25
Resolution	:	2.50 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:

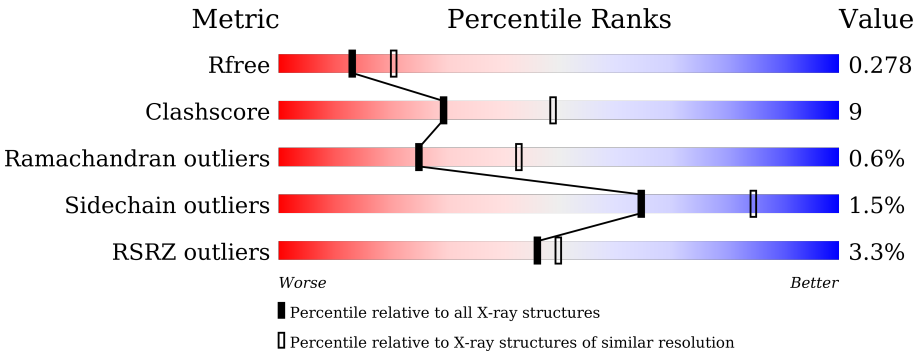
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.50 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	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

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 <=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	395	2% 71% 21% • 7%
1	B	395	4% 74% 15% 10%
1	C	395	3% 65% 21% • 13%
1	D	395	4% 66% 22% • 11%

2 Entry composition (i)

There are 2 unique types of molecules in this entry. The entry contains 10958 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 L-methionine gamma-lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	367	Total	C	N	O	S	0	0	0
			2817	1792	472	539	14			
1	B	355	Total	C	N	O	S	0	0	0
			2712	1727	455	518	12			
1	C	343	Total	C	N	O	S	0	0	0
			2611	1661	435	503	12			
1	D	350	Total	C	N	O	S	0	0	0
			2683	1710	451	508	14			

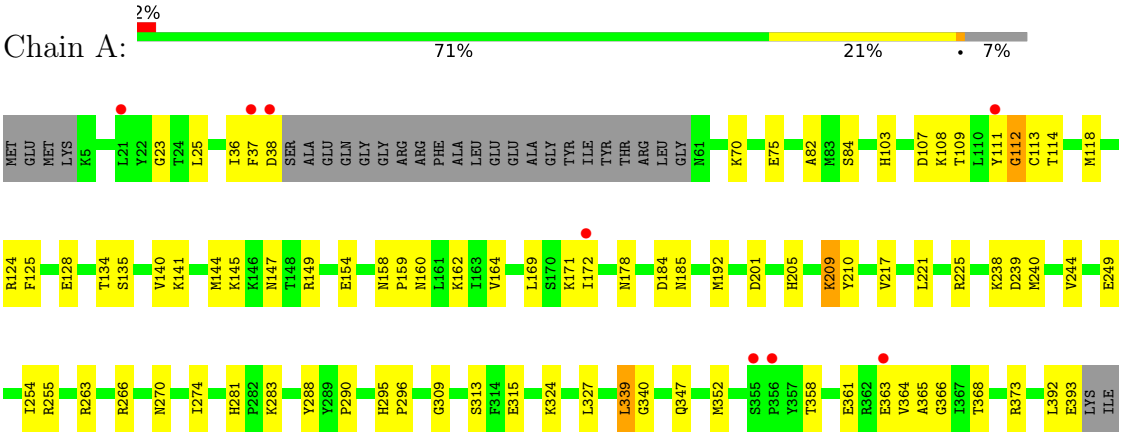
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	30	Total	O	0	0
			30	30		
2	B	35	Total	O	0	0
			35	35		
2	C	31	Total	O	0	0
			31	31		
2	D	39	Total	O	0	0
			39	39		

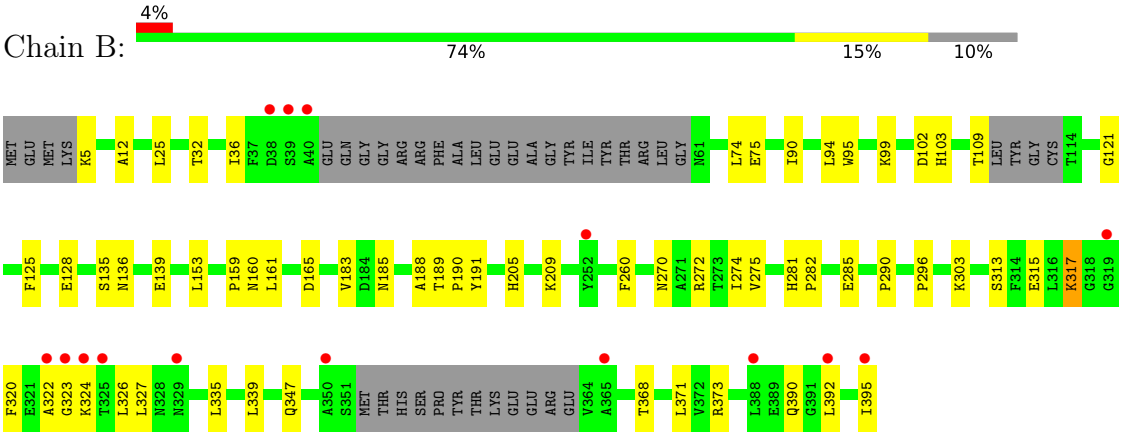
3 Residue-property plots (i)

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.

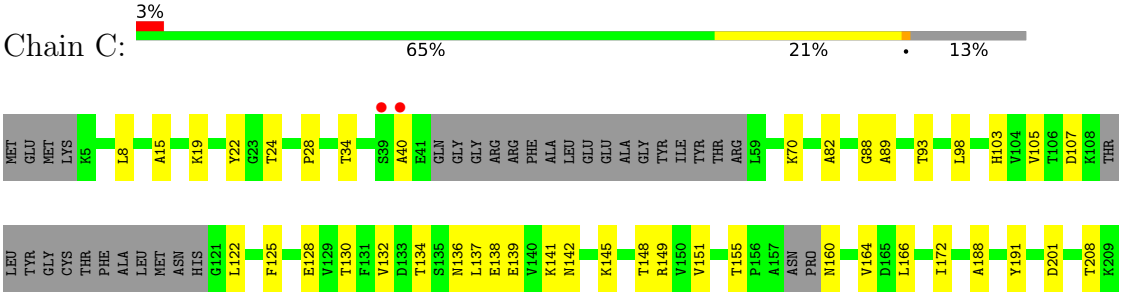
• Molecule 1: L-methionine gamma-lyase

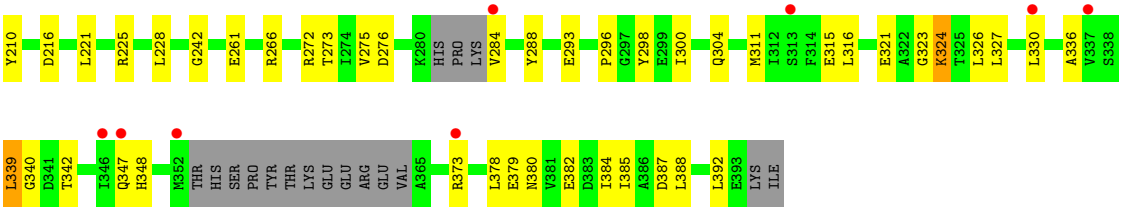


• Molecule 1: L-methionine gamma-lyase

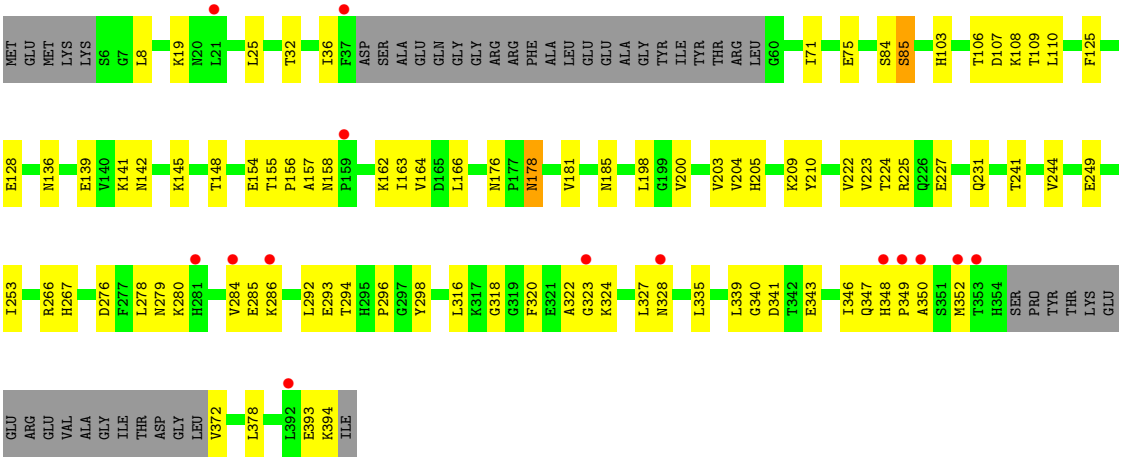


• Molecule 1: L-methionine gamma-lyase





• Molecule 1: L-methionine gamma-lyase



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	95.20Å 95.20Å 302.25Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	48.75 – 2.50 48.75 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.5 (48.75-2.50) 99.5 (48.75-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) >$ [1]	4.03 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.16_3549	Depositor
R, R_{free}	0.214 , 0.278 0.214 , 0.278	Depositor DCC
R_{free} test set	2751 reflections (4.92%)	wwPDB-VP
Wilson B-factor (Å^2)	47.7	Xtriage
Anisotropy	0.267	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 46.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	0.027 for -h,-k,l	Xtriage
F_o,F_c correlation	0.93	EDS
Total number of atoms	10958	wwPDB-VP
Average B, all atoms (Å^2)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.56% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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	0.42	0/2865	0.59	0/3881
1	B	0.43	0/2754	0.58	0/3726
1	C	0.42	0/2646	0.61	1/3575 (0.0%)
1	D	0.41	0/2728	0.56	0/3693
All	All	0.42	0/10993	0.58	1/14875 (0.0%)

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	188	ALA	C-N-CA	-5.76	107.30	121.70

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-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 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	2817	0	2862	53	0
1	B	2712	0	2771	41	0
1	C	2611	0	2664	65	0
1	D	2683	0	2735	69	0
2	A	30	0	0	1	0
2	B	35	0	0	1	0
2	C	31	0	0	5	0
2	D	39	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	10958	0	11032	207	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 9.

All (207) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:189:THR:HG22	1:B:191:TYR:H	1.38	0.88
1:A:141:LYS:HB2	1:A:172:ILE:HD11	1.61	0.81
1:B:322:ALA:O	1:B:395:ILE:HD13	1.84	0.77
1:B:32:THR:HG21	1:C:24:THR:HG21	1.65	0.77
1:C:382:GLU:H	1:C:382:GLU:CD	1.89	0.74
1:D:107:ASP:HB3	1:D:110:LEU:CD1	2.16	0.74
1:B:189:THR:HG23	1:B:260:PHE:HZ	1.52	0.74
1:D:176:ASN:OD1	1:D:178:ASN:ND2	2.21	0.73
1:C:201:ASP:OD1	1:C:225:ARG:NH1	2.23	0.71
1:C:142:ASN:OD1	2:C:401:HOH:O	2.08	0.71
1:C:89:ALA:O	1:C:93:THR:HG23	1.93	0.69
1:C:342:THR:HG22	1:C:378:LEU:HD12	1.73	0.69
1:D:279:ASN:HB3	1:D:280:LYS:HD3	1.76	0.68
1:D:284:VAL:HG12	1:D:316:LEU:HD23	1.76	0.66
1:C:40:ALA:HA	1:D:352:MET:HE2	1.78	0.65
1:B:188:ALA:O	1:B:189:THR:HB	1.96	0.65
1:B:161:LEU:HD23	1:B:290:PRO:HD3	1.78	0.65
1:C:40:ALA:H	1:D:328:ASN:HD21	1.45	0.64
1:B:368:THR:HG22	1:B:371:LEU:HB2	1.80	0.63
1:C:103:HIS:ND1	1:C:148:THR:OG1	2.28	0.63
1:C:330:LEU:HB3	1:C:387:ASP:OD2	1.97	0.63
1:B:189:THR:CG2	1:B:260:PHE:HZ	2.12	0.62
1:D:107:ASP:HB3	1:D:110:LEU:HD13	1.82	0.62
1:A:209:LYS:HG3	1:A:339:LEU:HG	1.82	0.62
1:C:34:THR:HG21	1:D:341:ASP:OD2	2.00	0.62
1:A:327:LEU:HD22	1:A:352:MET:HE1	1.82	0.61
1:D:348:HIS:HD2	1:D:372:VAL:HG22	1.64	0.61
1:C:155:THR:HB	1:C:164:VAL:HG22	1.83	0.60
1:D:209:LYS:HD2	1:D:339:LEU:HG	1.82	0.60
1:A:103:HIS:HD2	1:A:128:GLU:HB3	1.65	0.60
1:B:25:LEU:HD22	1:D:343:GLU:HG2	1.84	0.60
1:B:272:ARG:O	1:B:275:VAL:HG12	2.02	0.59
1:A:347:GLN:HG2	1:A:373:ARG:HB3	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:132:VAL:HB	1:C:139:GLU:HG2	1.84	0.59
1:A:25:LEU:HD11	1:D:36:ILE:HD11	1.83	0.59
1:B:136:ASN:HD22	1:B:139:GLU:HG3	1.67	0.59
1:D:162:LYS:HD3	1:D:163:ILE:H	1.67	0.58
1:D:348:HIS:CE1	1:D:350:ALA:HB3	2.38	0.58
1:A:270:ASN:O	1:A:274:ILE:HG13	2.04	0.58
1:D:103:HIS:HB3	1:D:148:THR:HA	1.86	0.58
1:D:142:ASN:ND2	2:D:405:HOH:O	2.36	0.57
1:B:5:LYS:N	2:B:406:HOH:O	2.37	0.56
1:A:327:LEU:HD22	1:A:352:MET:CE	2.36	0.56
1:A:108:LYS:NZ	1:A:109:THR:HG23	2.21	0.55
1:D:162:LYS:HZ1	1:D:296:PRO:HB2	1.70	0.55
1:A:201:ASP:OD1	1:A:225:ARG:NH1	2.39	0.55
1:C:347:GLN:NE2	1:C:373:ARG:HE	2.04	0.55
1:D:162:LYS:NZ	1:D:296:PRO:HB2	2.21	0.55
1:C:336:ALA:HB2	1:D:36:ILE:HD13	1.88	0.55
1:D:318:GLY:HA3	1:D:322:ALA:HB2	1.87	0.55
1:A:125:PHE:HB3	1:B:125:PHE:HB3	1.88	0.55
1:B:322:ALA:H	1:B:323:GLY:HA3	1.71	0.55
1:C:107:ASP:OD2	1:C:134:THR:OG1	2.24	0.55
1:B:136:ASN:ND2	1:B:139:GLU:HG3	2.22	0.54
1:A:315:GLU:OE2	1:A:368:THR:OG1	2.17	0.54
1:D:158:ASN:ND2	2:D:406:HOH:O	2.39	0.54
1:A:283:LYS:NZ	1:A:393:GLU:HA	2.23	0.54
1:A:217:VAL:HG13	1:A:249:GLU:HG2	1.89	0.54
1:D:136:ASN:HB3	1:D:139:GLU:HG2	1.89	0.54
1:D:203:VAL:HG12	1:D:223:VAL:HB	1.89	0.54
1:B:103:HIS:CD2	1:B:128:GLU:HG3	2.42	0.54
1:B:153:LEU:HB2	1:B:183:VAL:HG22	1.89	0.54
1:C:160:ASN:HB3	1:C:288:TYR:CD2	2.44	0.53
1:A:149:ARG:NH2	1:A:178:ASN:O	2.35	0.53
1:C:300:ILE:O	1:C:304:GLN:HG2	2.09	0.53
1:B:322:ALA:N	1:B:323:GLY:HA3	2.23	0.53
1:D:393:GLU:HG3	1:D:394:LYS:HD3	1.91	0.53
1:B:189:THR:HG23	1:B:260:PHE:CZ	2.40	0.52
1:D:103:HIS:HD2	1:D:128:GLU:HB3	1.73	0.52
1:A:36:ILE:CD1	1:D:19:LYS:HB2	2.40	0.52
1:A:36:ILE:HD11	1:D:25:LEU:HG	1.91	0.52
1:C:284:VAL:HG23	1:C:315:GLU:O	2.10	0.52
1:D:84:SER:HB3	1:D:244:VAL:O	2.10	0.52
1:C:326:LEU:HD13	1:C:392:LEU:HD23	1.91	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:105:VAL:CG1	1:C:151:VAL:HG22	2.40	0.52
1:D:107:ASP:OD1	1:D:108:LYS:N	2.43	0.52
1:A:290:PRO:HA	1:A:295:HIS:CD2	2.45	0.51
1:C:272:ARG:O	1:C:275:VAL:HG12	2.10	0.51
1:A:154:GLU:HG2	1:A:184:ASP:HB3	1.92	0.51
1:C:8:LEU:HD22	1:C:191:TYR:HD1	1.76	0.50
1:D:204:VAL:HG12	1:D:222:VAL:HG22	1.93	0.50
1:A:169:LEU:HA	1:A:172:ILE:HG22	1.93	0.50
1:C:276:ASP:OD2	2:C:402:HOH:O	2.19	0.50
1:D:278:LEU:O	1:D:284:VAL:HG21	2.12	0.50
1:D:349:PRO:HD3	1:D:372:VAL:HA	1.93	0.50
1:D:227:GLU:O	1:D:231:GLN:HG2	2.11	0.50
1:B:75:GLU:OE2	1:B:205:HIS:NE2	2.41	0.50
1:D:276:ASP:O	1:D:280:LYS:HG2	2.12	0.50
1:A:70:LYS:HE3	1:A:254:ILE:HG23	1.94	0.50
1:D:162:LYS:NZ	1:D:296:PRO:O	2.46	0.49
1:D:162:LYS:HA	1:D:162:LYS:HE2	1.94	0.49
1:B:159:PRO:HB3	1:B:373:ARG:CZ	2.42	0.49
1:B:189:THR:CG2	1:B:260:PHE:CZ	2.95	0.49
1:B:281:HIS:CD2	1:B:282:PRO:HD2	2.48	0.49
1:C:225:ARG:HH21	1:C:228:LEU:HD11	1.78	0.49
1:D:103:HIS:ND1	1:D:148:THR:OG1	2.29	0.49
1:C:15:ALA:O	1:C:70:LYS:NZ	2.43	0.49
1:C:105:VAL:HG13	1:C:151:VAL:HG22	1.95	0.48
1:C:323:GLY:HA3	1:C:348:HIS:CE1	2.48	0.48
1:B:99:LYS:NZ	1:B:102:ASP:OD1	2.46	0.48
1:A:160:ASN:HB3	1:A:288:TYR:CD2	2.49	0.48
1:B:103:HIS:HD2	1:B:128:GLU:HG3	1.77	0.48
1:C:216:ASP:HA	1:D:32:THR:HA	1.96	0.48
1:B:12:ALA:HB2	1:B:74:LEU:HD21	1.94	0.48
1:C:136:ASN:HB3	1:C:139:GLU:HB3	1.94	0.48
1:C:296:PRO:O	2:C:403:HOH:O	2.20	0.48
1:A:84:SER:HB3	1:A:244:VAL:O	2.13	0.48
1:D:156:PRO:HA	1:D:162:LYS:O	2.14	0.47
1:C:103:HIS:CD2	1:C:128:GLU:HG3	2.49	0.47
1:C:266:ARG:HD3	2:C:406:HOH:O	2.12	0.47
1:D:109:THR:O	1:D:110:LEU:HD12	2.14	0.47
1:C:141:LYS:HB2	1:C:172:ILE:HD13	1.95	0.47
1:C:103:HIS:HD2	1:C:128:GLU:HG3	1.80	0.47
1:A:107:ASP:OD2	1:A:134:THR:OG1	2.17	0.47
1:A:124:ARG:NH2	1:B:95:TRP:O	2.45	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:327:LEU:HA	1:C:330:LEU:CD1	2.45	0.47
1:C:125:PHE:HB3	1:D:125:PHE:HB3	1.96	0.47
1:A:75:GLU:OE2	1:A:205:HIS:NE2	2.42	0.46
1:C:82:ALA:HA	1:C:221:LEU:HD23	1.97	0.46
1:B:326:LEU:HD23	1:B:327:LEU:HD12	1.98	0.46
1:A:192:MET:HE1	1:A:309:GLY:HA2	1.97	0.46
1:A:112:GLY:O	1:A:113:CYS:HB2	2.14	0.46
1:C:98:LEU:HD23	1:C:149:ARG:HG2	1.97	0.46
1:C:242:GLY:O	1:D:85:SER:HB3	2.16	0.46
1:A:111:TYR:CD1	1:A:114:THR:HG23	2.51	0.45
1:A:185:ASN:HB3	1:A:205:HIS:CE1	2.51	0.45
1:C:107:ASP:HA	1:C:132:VAL:O	2.16	0.45
1:B:36:ILE:HD11	1:C:19:LYS:O	2.17	0.45
1:C:103:HIS:NE2	1:C:130:THR:HG22	2.30	0.45
1:B:165:ASP:OD2	1:B:303:LYS:NZ	2.47	0.45
1:B:327:LEU:HD23	1:B:335:LEU:HD12	1.98	0.45
1:D:249:GLU:O	1:D:253:ILE:HG13	2.17	0.45
1:C:293:GLU:HA	1:C:298:TYR:CG	2.51	0.45
1:C:284:VAL:HB	1:C:316:LEU:HD23	1.99	0.45
1:C:88:GLY:HA2	1:D:241:THR:O	2.17	0.44
1:C:321:GLU:O	1:C:324:LYS:HE3	2.16	0.44
1:D:285:GLU:HG3	1:D:286:LYS:HZ3	1.82	0.44
1:B:285:GLU:HA	1:B:317:LYS:HE3	1.99	0.44
1:D:155:THR:HB	1:D:164:VAL:HG22	2.00	0.44
1:A:25:LEU:CD1	1:D:36:ILE:HD11	2.48	0.44
1:A:288:TYR:HB2	1:A:313:SER:HB3	2.00	0.44
1:C:137:LEU:HD23	1:C:137:LEU:HA	1.82	0.44
1:C:273:THR:HG22	1:C:385:ILE:HD11	2.00	0.44
1:D:166:LEU:HD12	1:D:198:LEU:HD12	1.99	0.44
1:C:145:LYS:HB3	1:C:145:LYS:HE2	1.73	0.44
1:D:185:ASN:HB3	1:D:205:HIS:CE1	2.53	0.44
1:B:90:ILE:O	1:B:94:LEU:HG	2.18	0.44
1:D:267:HIS:CD2	1:D:378:LEU:HD21	2.53	0.44
1:D:346:ILE:HG13	1:D:347:GLN:N	2.32	0.44
1:A:358:THR:OG1	1:A:361:GLU:HB2	2.17	0.43
1:D:320:PHE:HE2	1:D:324:LYS:NZ	2.15	0.43
1:B:189:THR:CG2	1:B:191:TYR:H	2.21	0.43
1:C:22:TYR:CD1	1:C:28:PRO:HG2	2.53	0.43
1:D:106:THR:OG1	1:D:110:LEU:HD21	2.18	0.43
1:D:103:HIS:CE1	1:D:145:LYS:HD2	2.53	0.43
1:A:240:MET:CE	1:B:121:GLY:HA3	2.48	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:270:ASN:O	1:B:274:ILE:HG13	2.19	0.43
1:D:323:GLY:HA3	1:D:348:HIS:CD2	2.53	0.43
1:A:108:LYS:HZ3	1:A:109:THR:HG23	1.82	0.43
1:A:140:VAL:O	1:A:144:MET:HG3	2.18	0.43
1:C:272:ARG:HA	1:C:275:VAL:HG12	2.01	0.43
1:B:109:THR:O	1:B:109:THR:OG1	2.36	0.43
1:A:82:ALA:HA	1:A:221:LEU:HD23	2.00	0.43
1:A:255:ARG:HH21	1:C:216:ASP:CG	2.21	0.42
1:A:266:ARG:HH12	1:C:261:GLU:CD	2.23	0.42
1:B:190:PRO:HD3	1:B:205:HIS:CE1	2.54	0.42
1:D:224:THR:HG23	1:D:225:ARG:O	2.18	0.42
1:A:145:LYS:HE3	1:A:147:ASN:OD1	2.19	0.42
1:D:292:LEU:HB3	1:D:294:THR:HG22	2.01	0.42
1:D:75:GLU:OE2	1:D:205:HIS:NE2	2.47	0.42
1:A:263:ARG:HD2	2:A:412:HOH:O	2.18	0.42
1:D:162:LYS:HA	1:D:162:LYS:CE	2.49	0.42
1:A:365:ALA:HA	1:A:366:GLY:HA2	1.74	0.42
1:C:321:GLU:OE1	1:C:324:LYS:HE2	2.20	0.42
1:A:158:ASN:HA	1:A:159:PRO:HA	1.86	0.42
1:C:311:MET:SD	1:C:339:LEU:HB2	2.59	0.42
1:A:114:THR:O	1:A:118:MET:HG2	2.20	0.42
1:B:185:ASN:HB3	1:B:205:HIS:CE1	2.55	0.42
1:B:281:HIS:CE1	1:B:392:LEU:HD22	2.53	0.42
1:D:293:GLU:HA	1:D:298:TYR:CG	2.54	0.42
1:D:210:TYR:CE1	1:D:340:GLY:HA2	2.55	0.42
1:C:225:ARG:HB2	1:C:228:LEU:HD12	2.02	0.41
1:B:320:PHE:CE2	1:B:324:LYS:HD3	2.56	0.41
1:C:388:LEU:O	1:C:392:LEU:HG	2.20	0.41
1:A:172:ILE:HD13	1:A:172:ILE:HG21	1.81	0.41
1:A:281:HIS:CE1	1:A:392:LEU:HD13	2.55	0.41
1:C:315:GLU:OE1	2:C:404:HOH:O	2.22	0.41
1:A:37:PHE:HB3	1:A:38:ASP:H	1.64	0.41
1:A:164:VAL:HG13	1:A:169:LEU:CD1	2.50	0.41
1:D:181:VAL:HG13	1:D:200:VAL:HA	2.03	0.41
1:C:122:LEU:HA	1:C:122:LEU:HD23	1.79	0.41
1:C:128:GLU:H	1:C:128:GLU:HG2	1.76	0.41
1:C:210:TYR:CE1	1:C:340:GLY:HA2	2.56	0.41
1:D:178:ASN:HD22	1:D:178:ASN:H	1.69	0.41
1:A:23:GLY:O	1:D:36:ILE:N	2.48	0.41
1:A:364:VAL:HA	1:A:365:ALA:HA	1.88	0.41
1:A:210:TYR:CE1	1:A:340:GLY:HA2	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:135:SER:HB3	1:A:162:LYS:HE3	2.02	0.40
1:A:238:LYS:NZ	1:A:239:ASP:OD1	2.47	0.40
1:D:71:ILE:HD12	1:D:71:ILE:HA	1.90	0.40
1:D:154:GLU:O	1:D:157:ALA:HB2	2.21	0.40
1:D:327:LEU:O	1:D:335:LEU:HD11	2.21	0.40
1:C:166:LEU:H	1:C:304:GLN:HE22	1.68	0.40
1:D:109:THR:O	1:D:109:THR:OG1	2.38	0.40
1:C:321:GLU:HA	1:C:324:LYS:HG3	2.03	0.40
1:C:380:ASN:O	1:C:384:ILE:HG13	2.21	0.40
1:D:8:LEU:HD23	1:D:8:LEU:O	2.21	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	363/395 (92%)	344 (95%)	16 (4%)	3 (1%)	19	35
1	B	347/395 (88%)	331 (95%)	13 (4%)	3 (1%)	17	31
1	C	331/395 (84%)	311 (94%)	18 (5%)	2 (1%)	25	43
1	D	344/395 (87%)	326 (95%)	18 (5%)	0	100	100
All	All	1385/1580 (88%)	1312 (95%)	65 (5%)	8 (1%)	25	43

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	363	GLU
1	C	208	THR
1	B	317	LYS
1	B	209	LYS
1	B	296	PRO

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Mol	Chain	Res	Type
1	C	379	GLU
1	A	112	GLY
1	A	296	PRO

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	308/329 (94%)	304 (99%)	4 (1%)	69	87
1	B	296/329 (90%)	289 (98%)	7 (2%)	49	74
1	C	284/329 (86%)	281 (99%)	3 (1%)	73	89
1	D	293/329 (89%)	289 (99%)	4 (1%)	67	86
All	All	1181/1316 (90%)	1163 (98%)	18 (2%)	65	85

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	171	LYS
1	A	209	LYS
1	A	324	LYS
1	A	339	LEU
1	B	135	SER
1	B	160	ASN
1	B	313	SER
1	B	315	GLU
1	B	339	LEU
1	B	347	GLN
1	B	390	GLN
1	C	138	GLU
1	C	324	LYS
1	C	339	LEU
1	D	85	SER
1	D	141	LYS
1	D	178	ASN
1	D	266	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	231	GLN
1	B	281	HIS
1	C	304	GLN
1	C	347	GLN
1	D	178	ASN
1	D	328	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

There are no ligands in this entry.

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, 95^{th} 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	367/395 (92%)	-0.01	8 (2%)	62	65	28, 49, 75, 99	0
1	B	355/395 (89%)	0.08	15 (4%)	36	39	29, 46, 80, 104	0
1	C	343/395 (86%)	0.10	10 (2%)	51	55	31, 52, 77, 89	0
1	D	350/395 (88%)	0.14	14 (4%)	38	41	34, 52, 79, 102	0
All	All	1415/1580 (89%)	0.08	47 (3%)	46	50	28, 50, 79, 104	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	322	ALA	6.2
1	A	37	PHE	4.4
1	B	40	ALA	3.7
1	B	388	LEU	3.7
1	D	352	MET	3.6
1	A	356	PRO	3.6
1	C	284	VAL	3.2
1	A	172	ILE	3.2
1	B	392	LEU	3.2
1	C	40	ALA	3.2
1	D	353	THR	3.1
1	D	350	ALA	3.1
1	B	350	ALA	3.1
1	D	348	HIS	3.0
1	C	346	ILE	3.0
1	D	328	ASN	2.9
1	C	337	VAL	2.9
1	D	323	GLY	2.9
1	A	355	SER	2.9
1	B	395	ILE	2.9
1	B	323	GLY	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	365	ALA	2.7
1	B	39	SER	2.7
1	C	39	SER	2.6
1	C	373	ARG	2.6
1	D	284	VAL	2.5
1	A	21	LEU	2.5
1	D	281	HIS	2.5
1	C	347	GLN	2.5
1	C	330	LEU	2.4
1	A	38	ASP	2.4
1	B	324	LYS	2.3
1	A	111	TYR	2.2
1	C	313	SER	2.2
1	B	325	THR	2.2
1	B	329	ASN	2.2
1	A	363	GLU	2.1
1	B	38	ASP	2.1
1	D	37	PHE	2.1
1	D	21	LEU	2.1
1	C	352	MET	2.1
1	B	252	TYR	2.1
1	D	392	LEU	2.1
1	B	319	GLY	2.1
1	D	349	PRO	2.0
1	D	159	PRO	2.0
1	D	286	LYS	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

There are no ligands in this entry.

6.5 Other polymers (i)

There are no such residues in this entry.