Full wwPDB X-ray Structure Validation Report (i)

Nov 14, 2023 – 11:36 PM JST

PDB ID	:	6IGA
Title	:	Crystal structure of argininosuccinate lyase from Mycobacterium tuberculosis
Authors	:	Chen, X.B.; Liu, X.
Deposited on	:	2018-09-25
Resolution	:	2.78 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 (i) 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 (i)) were used in the production of this report:

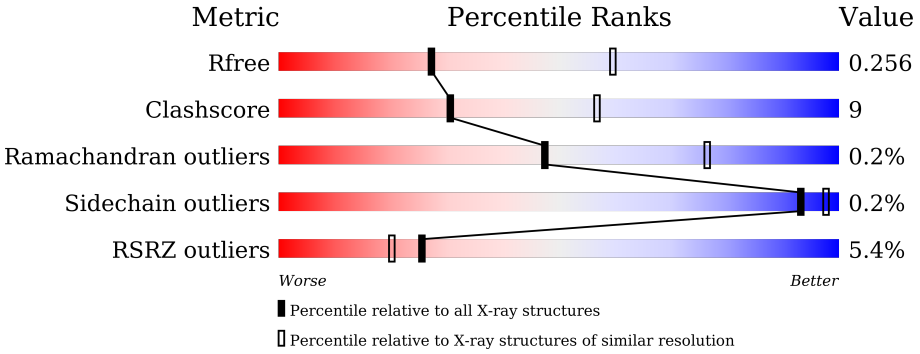
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, 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 2.78 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	4107 (2.80-2.76)
Clashscore	141614	4575 (2.80-2.76)
Ramachandran outliers	138981	4487 (2.80-2.76)
Sidechain outliers	138945	4489 (2.80-2.76)
RSRZ outliers	127900	4027 (2.80-2.76)

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	470	5% 78% 19% •
1	B	470	5% 76% 20% •
1	C	470	8% 77% 20% •
1	D	470	3% 81% 15% •

2 Entry composition (i)

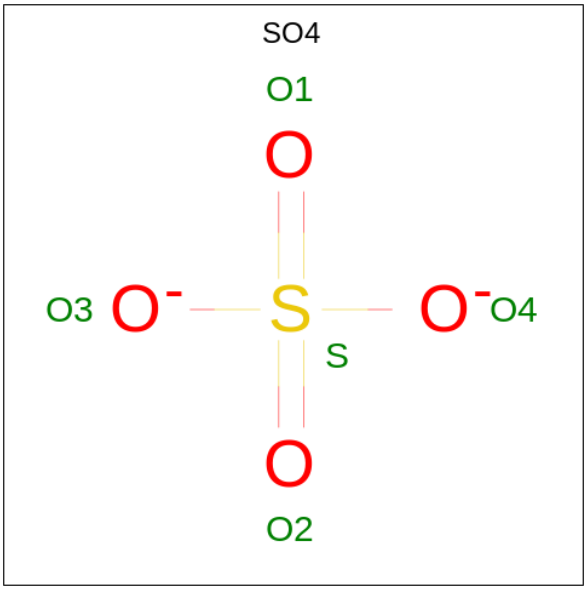
There are 2 unique types of molecules in this entry. The entry contains 13593 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Argininosuccinate lyase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	455	3397	2122	616	651	8	0	0	0
1	B	455	3397	2122	616	651	8	0	0	0
1	C	455	3397	2122	616	651	8	0	0	0
1	D	455	3397	2122	616	651	8	0	0	0

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



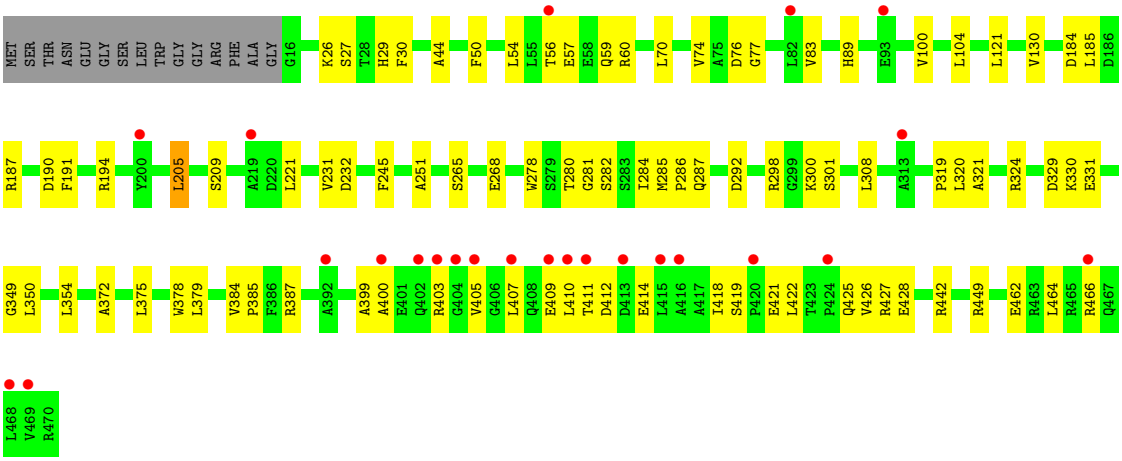
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
2	A	1	5	4	1	0	0

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.

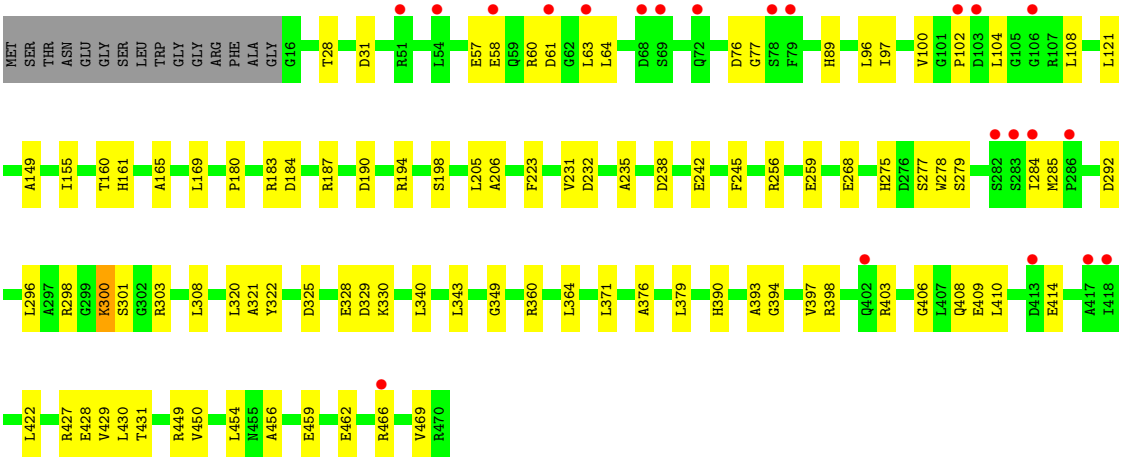
• Molecule 1: Argininosuccinate lyase

Chain A:



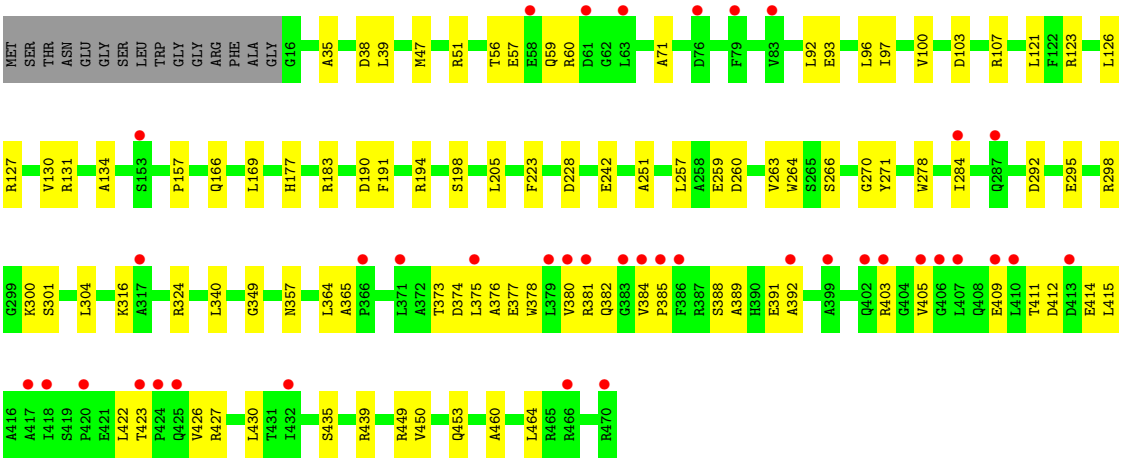
• Molecule 1: Argininosuccinate lyase

Chain B:



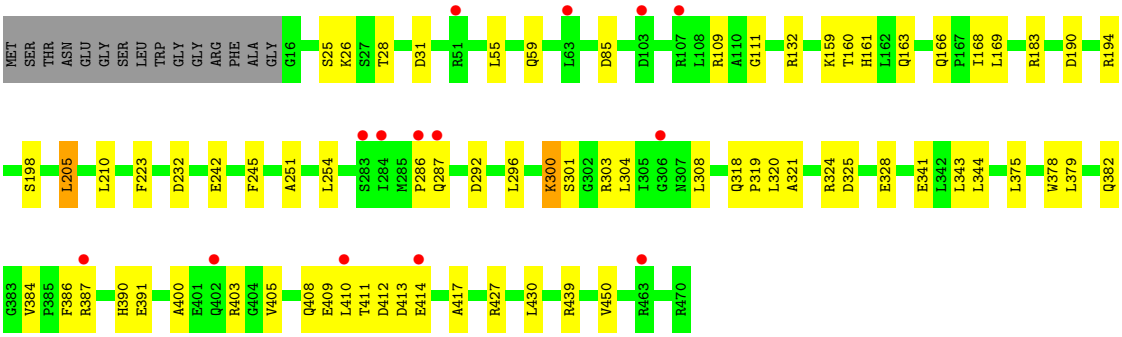
• Molecule 1: Argininosuccinate lyase

Chain C:



- Molecule 1: Argininosuccinate lyase

Chain D:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α, β, γ	129.16Å 129.16Å 200.14Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	45.67 – 2.78 45.67 – 2.78	Depositor EDS
% Data completeness (in resolution range)	88.9 (45.67-2.78) 88.9 (45.67-2.78)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.42 (at 2.77Å)	Xtriage
Refinement program	PHENIX 1.8.4_1496	Depositor
R, R_{free}	0.219 , 0.255 0.220 , 0.256	Depositor DCC
R_{free} test set	2170 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å^2)	43.6	Xtriage
Anisotropy	0.266	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 24.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.033 for -h,-k,l	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	13593	wwPDB-VP
Average B, all atoms (Å^2)	48.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.82% 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 (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: SO4

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\#\|Z\| >5$	RMSZ	$\#\|Z\| >5$
1	A	0.40	0/3455	0.50	0/4697
1	B	0.39	0/3455	0.46	0/4697
1	C	0.32	0/3455	0.48	0/4697
1	D	0.29	0/3455	0.45	0/4697
All	All	0.35	0/13820	0.47	0/18788

There are no bond length outliers.

There are no bond angle outliers.

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	3397	0	3388	70	0
1	B	3397	0	3388	71	0
1	C	3397	0	3388	84	0
1	D	3397	0	3388	60	0
2	A	5	0	0	0	0
All	All	13593	0	13552	250	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 (250) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:405:VAL:HB	1:C:409:GLU:HB2	1.61	0.83
1:D:386:PHE:CE1	1:D:390:HIS:CE1	2.69	0.80
1:A:287:GLN:OE1	1:A:287:GLN:N	2.15	0.76
1:D:160:THR:HG22	1:D:161:HIS:CD2	2.21	0.76
1:C:126:LEU:O	1:C:130:VAL:HG23	1.85	0.76
1:C:376:ALA:O	1:C:380:VAL:HG23	1.87	0.75
1:A:387:ARG:NH2	1:C:103:ASP:OD1	2.20	0.74
1:D:109:ARG:HG2	1:D:109:ARG:HH11	1.52	0.74
1:C:388:SER:O	1:C:391:GLU:HB2	1.88	0.73
1:A:190:ASP:OD2	1:C:194:ARG:NH2	2.22	0.72
1:B:194:ARG:NH2	1:D:190:ASP:OD2	2.23	0.72
1:C:388:SER:HA	1:C:391:GLU:HB2	1.72	0.71
1:A:442:ARG:NH2	1:C:228:ASP:OD1	2.22	0.71
1:D:387:ARG:O	1:D:391:GLU:HG2	1.91	0.70
1:B:232:ASP:OD2	1:D:183:ARG:NH2	2.25	0.68
1:A:324:ARG:NH2	1:B:292:ASP:OD2	2.26	0.68
1:B:183:ARG:NH2	1:D:232:ASP:OD2	2.27	0.68
1:C:460:ALA:O	1:C:464:LEU:HG	1.94	0.68
1:C:166:GLN:HG2	1:C:439:ARG:NH2	2.10	0.67
1:D:25:SER:OG	1:D:324:ARG:NH2	2.28	0.67
1:A:268:GLU:OE2	1:D:159:LYS:NZ	2.27	0.67
1:C:324:ARG:NH1	1:D:292:ASP:OD2	2.28	0.67
1:A:26:LYS:NZ	1:A:83:VAL:O	2.24	0.66
1:A:286:PRO:HD2	1:A:287:GLN:OE1	1.95	0.66
1:B:403:ARG:NH2	1:B:414:GLU:OE2	2.29	0.66
1:A:185:LEU:HB2	1:A:464:LEU:HD13	1.79	0.65
1:C:415:LEU:O	1:C:422:LEU:HD12	1.97	0.65
1:B:300:LYS:HD2	1:B:343:LEU:HD11	1.79	0.65
1:D:411:THR:OG1	1:D:414:GLU:HG3	1.98	0.64
1:C:166:GLN:HG2	1:C:439:ARG:HH21	1.62	0.64
1:B:194:ARG:HD3	1:D:183:ARG:HG2	1.80	0.64
1:C:375:LEU:HD12	1:C:430:LEU:HD21	1.80	0.63
1:A:300:LYS:NZ	1:B:328:GLU:OE2	2.32	0.62
1:B:371:LEU:HD13	1:B:430:LEU:O	1.98	0.62
1:C:93:GLU:HG3	1:C:97:ILE:HD13	1.81	0.62
1:B:60:ARG:O	1:B:64:LEU:HG	1.99	0.62
1:B:394:GLY:O	1:B:398:ARG:HG3	1.98	0.62
1:B:410:LEU:O	1:B:427:ARG:NH1	2.33	0.61
1:A:100:VAL:HB	1:A:104:LEU:HD23	1.81	0.61
1:B:61:ASP:HA	1:B:64:LEU:HB2	1.81	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:50:PHE:HE1	1:A:57:GLU:HG2	1.66	0.61
1:B:190:ASP:OD2	1:D:194:ARG:NH2	2.35	0.60
1:A:462:GLU:O	1:A:466:ARG:HG3	2.01	0.59
1:B:97:ILE:HG23	1:B:102:PRO:HA	1.84	0.59
1:A:232:ASP:OD2	1:C:183:ARG:NH2	2.35	0.59
1:D:296:LEU:O	1:D:300:LYS:HD3	2.02	0.59
1:A:287:GLN:H	1:A:287:GLN:CD	2.05	0.59
1:A:384:VAL:CG2	1:A:421:GLU:HG2	2.32	0.59
1:C:377:GLU:O	1:C:381:ARG:HG3	2.02	0.59
1:C:388:SER:O	1:C:392:ALA:N	2.29	0.58
1:B:408:GLN:NE2	1:B:431:THR:HG23	2.18	0.58
1:A:282:SER:OG	1:A:284:ILE:HG13	2.04	0.57
1:C:403:ARG:NH1	1:C:405:VAL:HG11	2.18	0.57
1:C:177:HIS:NE2	1:C:260:ASP:OD2	2.31	0.57
1:A:411:THR:HG22	1:A:412:ASP:N	2.20	0.56
1:C:378:TRP:HD1	1:C:381:ARG:HH11	1.52	0.56
1:B:322:TYR:HB3	1:D:160:THR:HG23	1.88	0.56
1:C:378:TRP:HD1	1:C:381:ARG:NH1	2.04	0.56
1:A:89:HIS:HE1	1:A:121:LEU:HD12	1.71	0.56
1:C:388:SER:CA	1:C:391:GLU:HB2	2.35	0.56
1:C:423:THR:O	1:C:426:VAL:HG12	2.06	0.56
1:C:378:TRP:CE3	1:C:426:VAL:HG23	2.41	0.56
1:B:393:ALA:O	1:B:397:VAL:HG23	2.06	0.55
1:D:378:TRP:CZ3	1:D:382:GLN:HG3	2.41	0.55
1:A:419:SER:HB3	1:A:422:LEU:HD13	1.88	0.55
1:C:300:LYS:NZ	1:D:328:GLU:OE2	2.22	0.55
1:C:166:GLN:HE21	1:C:373:THR:HG21	1.70	0.55
1:A:378:TRP:CE3	1:A:426:VAL:HG23	2.42	0.54
1:B:275:HIS:O	1:B:279:SER:OG	2.25	0.54
1:D:400:ALA:CB	1:D:410:LEU:HD21	2.38	0.54
1:A:44:ALA:HB2	1:A:221:LEU:HD21	1.89	0.54
1:C:130:VAL:HG21	1:C:191:PHE:CE2	2.44	0.53
1:C:382:GLN:HA	1:C:382:GLN:OE1	2.06	0.53
1:A:384:VAL:HG21	1:A:421:GLU:HG2	1.90	0.53
1:C:194:ARG:NH1	1:C:242:GLU:OE2	2.41	0.53
1:C:411:THR:O	1:C:414:GLU:HB2	2.08	0.53
1:C:374:ASP:OD2	1:C:435:SER:OG	2.26	0.53
1:C:59:GLN:HE22	1:C:107:ARG:HH11	1.55	0.53
1:D:412:ASP:OD1	1:D:427:ARG:NH2	2.27	0.53
1:D:408:GLN:HG2	1:D:409:GLU:N	2.24	0.52
1:C:38:ASP:HB2	1:C:121:LEU:HD21	1.90	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:379:LEU:HD22	1:D:384:VAL:HG21	1.91	0.52
1:D:400:ALA:HB2	1:D:410:LEU:HD21	1.92	0.52
1:A:265:SER:O	1:D:163:GLN:NE2	2.36	0.52
1:D:405:VAL:HG22	1:D:409:GLU:HB2	1.91	0.52
1:A:324:ARG:HD2	1:B:292:ASP:HB3	1.90	0.52
1:B:428:GLU:HG3	1:B:429:VAL:HG13	1.90	0.52
1:C:384:VAL:CG1	1:C:385:PRO:HD2	2.40	0.52
1:D:251:ALA:HB2	1:D:304:LEU:HD22	1.92	0.52
1:C:384:VAL:CG1	1:C:385:PRO:CD	2.87	0.52
1:D:251:ALA:HB1	1:D:301:SER:HA	1.91	0.51
1:A:286:PRO:CD	1:A:287:GLN:OE1	2.58	0.51
1:B:169:LEU:HD12	1:B:450:VAL:HG21	1.91	0.51
1:A:379:LEU:HD11	1:A:422:LEU:HD11	1.91	0.51
1:D:296:LEU:O	1:D:300:LYS:HB2	2.11	0.51
1:A:278:TRP:CZ3	1:A:349:GLY:HA3	2.45	0.51
1:B:76:ASP:OD1	1:B:77:GLY:N	2.44	0.51
1:D:400:ALA:HA	1:D:410:LEU:HD21	1.93	0.51
1:A:209:SER:O	1:C:381:ARG:HD3	2.11	0.51
1:A:76:ASP:OD1	1:A:77:GLY:N	2.44	0.51
1:B:298:ARG:O	1:B:301:SER:OG	2.28	0.51
1:D:378:TRP:CH2	1:D:382:GLN:HG3	2.46	0.51
1:A:54:LEU:HD11	1:C:381:ARG:O	2.11	0.50
1:C:264:TRP:CD1	1:C:271:TYR:HD2	2.29	0.50
1:C:388:SER:C	1:C:391:GLU:HB2	2.31	0.50
1:A:285:MET:HG2	1:A:287:GLN:HE22	1.76	0.50
1:B:360:ARG:HE	1:B:364:LEU:HD11	1.77	0.50
1:D:245:PHE:HA	1:D:308:LEU:HD22	1.94	0.50
1:B:379:LEU:HD21	1:B:422:LEU:HD23	1.94	0.50
1:D:386:PHE:CZ	1:D:390:HIS:CE1	3.00	0.50
1:C:57:GLU:CD	1:C:60:ARG:HE	2.13	0.50
1:B:245:PHE:HA	1:B:308:LEU:HD22	1.92	0.49
1:D:403:ARG:NH1	1:D:411:THR:HG23	2.26	0.49
1:B:194:ARG:NH1	1:B:242:GLU:OE2	2.44	0.49
1:B:284:ILE:HB	1:B:285:MET:HE2	1.93	0.49
1:B:296:LEU:O	1:B:300:LYS:HB2	2.11	0.49
1:A:251:ALA:HB1	1:A:301:SER:HA	1.94	0.49
1:B:278:TRP:CH2	1:B:349:GLY:HA3	2.46	0.49
1:D:194:ARG:NH1	1:D:242:GLU:OE2	2.46	0.49
1:D:109:ARG:HG2	1:D:109:ARG:NH1	2.23	0.49
1:C:378:TRP:HA	1:C:381:ARG:CZ	2.42	0.49
1:B:409:GLU:O	1:B:410:LEU:HD12	2.13	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:206:ALA:HB2	1:D:168:ILE:HG22	1.94	0.49
1:B:456:ALA:O	1:B:459:GLU:HG2	2.13	0.48
1:C:384:VAL:HG12	1:C:385:PRO:HD2	1.95	0.48
1:B:376:ALA:HB1	1:B:390:HIS:CD2	2.48	0.48
1:B:275:HIS:ND1	1:B:277:SER:HB2	2.29	0.48
1:B:180:PRO:HG2	1:B:256:ARG:HH12	1.78	0.48
1:D:413:ASP:O	1:D:417:ALA:N	2.42	0.48
1:A:403:ARG:HB2	1:A:405:VAL:HG13	1.95	0.48
1:C:56:THR:OG1	1:C:59:GLN:HG3	2.13	0.48
1:C:130:VAL:HG21	1:C:191:PHE:HE2	1.78	0.48
1:D:300:LYS:CB	1:D:343:LEU:HD11	2.44	0.48
1:B:96:LEU:O	1:B:100:VAL:HG22	2.13	0.47
1:C:378:TRP:CD1	1:C:381:ARG:NH1	2.82	0.47
1:C:123:ARG:O	1:C:127:ARG:HG3	2.13	0.47
1:C:384:VAL:HG12	1:C:385:PRO:CD	2.44	0.47
1:A:449:ARG:HA	1:A:449:ARG:HD3	1.70	0.47
1:C:298:ARG:O	1:C:301:SER:OG	2.23	0.47
1:A:50:PHE:CD1	1:A:60:ARG:HD3	2.50	0.47
1:A:399:ALA:O	1:A:400:ALA:C	2.53	0.47
1:B:408:GLN:NE2	1:B:431:THR:CG2	2.78	0.47
1:C:39:LEU:HD13	1:C:71:ALA:HA	1.97	0.47
1:C:47:MET:O	1:C:51:ARG:HG3	2.15	0.47
1:A:56:THR:O	1:A:60:ARG:N	2.43	0.47
1:B:149:ALA:HB2	1:B:454:LEU:HD21	1.97	0.47
1:B:406:GLY:O	1:B:410:LEU:HD13	2.15	0.46
1:A:26:LYS:HZ1	1:A:83:VAL:HA	1.80	0.46
1:A:278:TRP:CH2	1:A:349:GLY:HA3	2.49	0.46
1:A:285:MET:HG2	1:A:287:GLN:NE2	2.30	0.46
1:C:96:LEU:O	1:C:100:VAL:HG22	2.16	0.46
1:A:414:GLU:O	1:A:418:ILE:HG12	2.16	0.46
1:B:462:GLU:O	1:B:466:ARG:HG3	2.16	0.46
1:A:130:VAL:HG21	1:A:191:PHE:CE2	2.51	0.46
1:B:238:ASP:OD2	1:D:183:ARG:HD2	2.17	0.45
1:B:390:HIS:ND1	1:C:284:ILE:HD11	2.31	0.45
1:C:198:SER:HB2	1:C:223:PHE:CD2	2.51	0.45
1:A:56:THR:O	1:A:59:GLN:HB2	2.16	0.45
1:A:378:TRP:CD2	1:A:426:VAL:HG23	2.51	0.45
1:B:449:ARG:HA	1:B:449:ARG:HD3	1.80	0.45
1:D:132:ARG:NE	1:D:341:GLU:OE2	2.40	0.45
1:B:63:LEU:HD13	1:B:108:LEU:HD13	1.97	0.45
1:C:198:SER:HB2	1:C:223:PHE:CG	2.51	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:184:ASP:OD1	1:A:187:ARG:NH1	2.49	0.45
1:D:28:THR:HA	1:D:31:ASP:OD1	2.16	0.45
1:D:386:PHE:CE1	1:D:390:HIS:HE1	2.31	0.45
1:C:157:PRO:HB3	1:C:365:ALA:HA	1.98	0.45
1:C:378:TRP:CD2	1:C:426:VAL:HG23	2.52	0.45
1:D:286:PRO:HB2	1:D:287:GLN:NE2	2.31	0.45
1:A:411:THR:HG22	1:A:412:ASP:H	1.81	0.45
1:C:131:ARG:O	1:C:134:ALA:HB3	2.17	0.45
1:A:245:PHE:HA	1:A:308:LEU:HD22	1.99	0.45
1:B:300:LYS:CD	1:B:343:LEU:HD11	2.45	0.45
1:A:194:ARG:NH2	1:C:190:ASP:OD2	2.50	0.45
1:B:198:SER:HB2	1:B:223:PHE:CG	2.52	0.45
1:D:166:GLN:HG2	1:D:439:ARG:NH2	2.32	0.45
1:A:298:ARG:O	1:C:316:LYS:HE2	2.16	0.44
1:C:169:LEU:HD12	1:C:450:VAL:HG21	1.98	0.44
1:B:100:VAL:HB	1:B:104:LEU:HD22	1.99	0.44
1:B:198:SER:HB2	1:B:223:PHE:CD2	2.51	0.44
1:C:270:GLY:O	1:C:357:ASN:ND2	2.50	0.44
1:A:324:ARG:HD3	1:B:296:LEU:HG	1.99	0.44
1:A:29:HIS:CE1	1:A:30:PHE:HD2	2.35	0.44
1:D:300:LYS:O	1:D:303:ARG:HB3	2.17	0.44
1:C:384:VAL:HG12	1:C:385:PRO:N	2.33	0.44
1:C:389:ALA:O	1:C:392:ALA:HB3	2.17	0.44
1:B:89:HIS:HE1	1:B:121:LEU:HD12	1.83	0.43
1:C:427:ARG:O	1:C:430:LEU:HB2	2.18	0.43
1:A:425:GLN:O	1:A:428:GLU:HB2	2.19	0.43
1:C:295:GLU:HB3	1:D:325:ASP:OD1	2.18	0.43
1:A:331:GLU:OE2	1:B:303:ARG:NH1	2.46	0.43
1:A:385:PRO:HB3	1:C:107:ARG:HG2	1.99	0.43
1:B:184:ASP:OD1	1:B:187:ARG:NH1	2.52	0.43
1:D:375:LEU:HD12	1:D:430:LEU:HD21	2.00	0.43
1:C:59:GLN:HE22	1:C:107:ARG:NH1	2.16	0.43
1:C:292:ASP:OD2	1:D:324:ARG:NH1	2.51	0.43
1:D:111:GLY:HA3	1:D:210:LEU:HD22	2.00	0.43
1:C:271:TYR:HE1	1:C:364:LEU:HD12	1.83	0.43
1:D:400:ALA:CA	1:D:410:LEU:HD21	2.48	0.43
1:C:449:ARG:HD3	1:C:449:ARG:HA	1.66	0.43
1:A:280:THR:CG2	1:A:281:GLY:N	2.81	0.43
1:D:405:VAL:HG22	1:D:409:GLU:CB	2.48	0.43
1:A:298:ARG:HB3	1:B:325:ASP:OD1	2.19	0.42
1:B:165:ALA:HB3	1:D:205:LEU:HD21	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:160:THR:HG22	1:B:161:HIS:ND1	2.34	0.42
1:C:388:SER:O	1:C:391:GLU:N	2.52	0.42
1:C:340:LEU:HD23	1:C:340:LEU:HA	1.90	0.42
1:C:388:SER:O	1:C:391:GLU:CB	2.64	0.42
1:B:231:VAL:HG13	1:B:321:ALA:HB2	2.02	0.42
1:B:268:GLU:OE2	1:C:271:TYR:OH	2.36	0.42
1:D:198:SER:HB2	1:D:223:PHE:CG	2.54	0.42
1:A:372:ALA:O	1:A:375:LEU:HB3	2.18	0.42
1:A:410:LEU:O	1:A:427:ARG:NH2	2.53	0.42
1:B:155:ILE:HG12	1:B:169:LEU:HD23	2.02	0.42
1:B:235:ALA:HB2	1:B:321:ALA:HB1	2.02	0.42
1:C:278:TRP:CZ3	1:C:349:GLY:HA3	2.54	0.42
1:D:169:LEU:HD12	1:D:450:VAL:HG21	2.00	0.42
1:A:320:LEU:HA	1:A:321:ALA:HA	1.76	0.42
1:A:205:LEU:HD12	1:A:231:VAL:HG22	2.02	0.42
1:B:259:GLU:HG2	1:D:319:PRO:HD3	2.01	0.42
1:B:278:TRP:CZ3	1:B:349:GLY:HA3	2.55	0.42
1:A:319:PRO:HD3	1:C:259:GLU:HG2	2.01	0.41
1:C:35:ALA:HA	1:C:92:LEU:HD11	2.02	0.41
1:C:251:ALA:HB2	1:C:304:LEU:HD22	2.01	0.41
1:B:57:GLU:O	1:B:58:GLU:C	2.54	0.41
1:D:55:LEU:HD22	1:D:59:GLN:HB3	2.01	0.41
1:D:320:LEU:HA	1:D:321:ALA:HA	1.81	0.41
1:A:27:SER:HB2	1:A:30:PHE:CZ	2.55	0.41
1:A:70:LEU:O	1:A:74:VAL:HG23	2.21	0.41
1:B:340:LEU:HD23	1:B:340:LEU:HA	1.91	0.41
1:A:280:THR:HB	1:A:292:ASP:OD2	2.21	0.41
1:A:329:ASP:OD1	1:A:330:LYS:N	2.53	0.41
1:B:245:PHE:CZ	1:D:242:GLU:HA	2.55	0.41
1:B:379:LEU:HD23	1:B:379:LEU:HA	1.91	0.41
1:B:28:THR:HA	1:B:31:ASP:OD1	2.20	0.41
1:C:450:VAL:HA	1:C:453:GLN:HG2	2.03	0.41
1:D:318:GLN:NE2	1:D:325:ASP:OD2	2.42	0.41
1:A:282:SER:OG	1:A:284:ILE:CG1	2.68	0.41
1:A:379:LEU:HD23	1:A:379:LEU:HA	1.93	0.41
1:A:400:ALA:CB	1:A:407:LEU:HD23	2.51	0.41
1:B:320:LEU:HA	1:B:321:ALA:HA	1.65	0.41
1:C:263:VAL:O	1:C:266:SER:OG	2.32	0.41
1:D:26:LYS:NZ	1:D:85:ASP:O	2.53	0.41
1:D:254:LEU:HD11	1:D:344:LEU:HD23	2.03	0.41
1:C:177:HIS:CE1	1:C:257:LEU:HA	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:384:VAL:HG13	1:C:385:PRO:CD	2.51	0.40
1:C:412:ASP:OD1	1:C:427:ARG:NH1	2.50	0.40
1:A:350:LEU:O	1:A:354:LEU:HG	2.20	0.40
1:B:466:ARG:O	1:B:469:VAL:HG22	2.21	0.40
1:B:329:ASP:OD1	1:B:330:LYS:HG3	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	453/470 (96%)	441 (97%)	11 (2%)	1 (0%)	47	76
1	B	453/470 (96%)	443 (98%)	9 (2%)	1 (0%)	47	76
1	C	453/470 (96%)	438 (97%)	14 (3%)	1 (0%)	47	76
1	D	453/470 (96%)	436 (96%)	16 (4%)	1 (0%)	47	76
All	All	1812/1880 (96%)	1758 (97%)	50 (3%)	4 (0%)	47	76

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	205	LEU
1	B	205	LEU
1	C	205	LEU
1	D	205	LEU

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	342/352 (97%)	341 (100%)	1 (0%)	92	97
1	B	342/352 (97%)	341 (100%)	1 (0%)	92	97
1	C	342/352 (97%)	342 (100%)	0	100	100
1	D	342/352 (97%)	341 (100%)	1 (0%)	92	97
All	All	1368/1408 (97%)	1365 (100%)	3 (0%)	93	98

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	409	GLU
1	B	300	LYS
1	D	300	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	C	390	HIS
1	D	161	HIS
1	D	390	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	SO4	A	501	-	4,4,4	0.32	0	6,6,6	0.05	0

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 (i)

6.1 Protein, DNA and RNA chains (i)

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	455/470 (96%)	0.34	24 (5%)	26	21	26, 48, 90, 123	0
1	B	455/470 (96%)	0.22	22 (4%)	30	24	25, 41, 79, 111	0
1	C	455/470 (96%)	0.43	39 (8%)	10	7	27, 49, 104, 131	0
1	D	455/470 (96%)	0.08	14 (3%)	49	44	22, 39, 71, 102	0
All	All	1820/1880 (96%)	0.27	99 (5%)	25	20	22, 44, 87, 131	0

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	404	GLY	5.9
1	C	405	VAL	5.9
1	A	400	ALA	5.8
1	C	470	ARG	5.4
1	B	284	ILE	5.2
1	B	102	PRO	5.1
1	C	407	LEU	5.0
1	A	410	LEU	5.0
1	A	413	ASP	4.9
1	B	58	GLU	4.7
1	A	402	GLN	4.7
1	A	411	THR	4.6
1	C	403	ARG	4.6
1	B	61	ASP	4.6
1	B	283	SER	4.6
1	C	375	LEU	4.5
1	C	413	ASP	4.3
1	A	409	GLU	4.1
1	B	417	ALA	3.9
1	D	284	ILE	3.9
1	C	424	PRO	3.9

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Mol	Chain	Res	Type	RSRZ
1	C	399	ALA	3.8
1	D	286	PRO	3.8
1	C	380	VAL	3.7
1	B	103	ASP	3.6
1	B	72	GLN	3.6
1	C	153	SER	3.5
1	B	51	ARG	3.5
1	A	405	VAL	3.5
1	A	407	LEU	3.4
1	C	410	LEU	3.4
1	A	466	ARG	3.4
1	B	466	ARG	3.2
1	B	63	LEU	3.1
1	B	402	GLN	3.1
1	B	286	PRO	3.1
1	C	425	GLN	3.0
1	B	68	ASP	3.0
1	D	402	GLN	3.0
1	C	420	PRO	3.0
1	C	381	ARG	3.0
1	C	466	ARG	2.9
1	C	386	PHE	2.9
1	A	415	LEU	2.8
1	C	76	ASP	2.8
1	A	420	PRO	2.7
1	D	287	GLN	2.7
1	A	403	ARG	2.7
1	A	424	PRO	2.7
1	C	402	GLN	2.7
1	C	284	ILE	2.7
1	A	56	THR	2.7
1	B	106	GLY	2.7
1	B	418	ILE	2.7
1	C	287	GLN	2.6
1	C	392	ALA	2.6
1	B	54	LEU	2.6
1	C	418	ILE	2.6
1	C	379	LEU	2.5
1	B	78	SER	2.5
1	A	93	GLU	2.5
1	D	410	LEU	2.5
1	C	58	GLU	2.5

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Mol	Chain	Res	Type	RSRZ
1	C	417	ALA	2.5
1	C	409	GLU	2.4
1	C	83	VAL	2.4
1	D	283	SER	2.4
1	C	385	PRO	2.4
1	A	219	ALA	2.4
1	D	463	ARG	2.4
1	A	82	LEU	2.3
1	C	63	LEU	2.3
1	D	103	ASP	2.3
1	C	317	ALA	2.3
1	C	366	PRO	2.3
1	C	432	ILE	2.3
1	A	468	LEU	2.3
1	A	392	ALA	2.3
1	B	282	SER	2.2
1	D	107	ARG	2.2
1	A	416	ALA	2.2
1	D	387	ARG	2.2
1	C	423	THR	2.2
1	D	306	GLY	2.2
1	C	371	LEU	2.1
1	A	313	ALA	2.1
1	D	414	GLU	2.1
1	D	51	ARG	2.1
1	B	69	SER	2.1
1	C	61	ASP	2.1
1	C	406	GLY	2.1
1	A	469	VAL	2.1
1	C	79	PHE	2.1
1	C	384	VAL	2.0
1	C	383	GLY	2.0
1	D	63	LEU	2.0
1	A	200	TYR	2.0
1	B	413	ASP	2.0
1	B	79	PHE	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(Å^2)	Q<0.9
2	SO4	A	501	5/5	0.87	0.34	33,33,33,33	5

6.5 Other polymers (i)

There are no such residues in this entry.