Full wwPDB X-ray Structure Validation Report (i)

Jun 19, 2024 – 08:00 AM EDT

PDB ID	:	4NNP
Title	:	Crystal Structure of Apo Manganese ABC transporter MntC from Staphylococcus aureus bound to an antagonistic fab fragment
Authors	:	Rouge, L.; Sudhamsu, J.
Deposited on	:	2013-11-18
Resolution	:	2.69 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 (1)) were used in the production of this report:

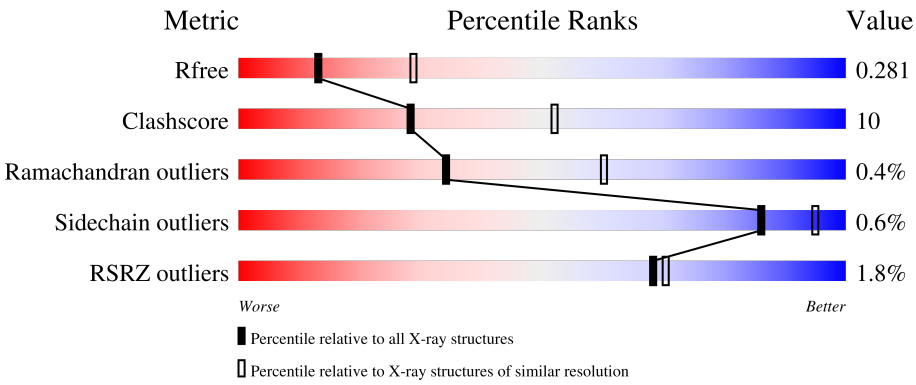
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation 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 2.69 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	2808 (2.70-2.70)
Clashscore	141614	3122 (2.70-2.70)
Ramachandran outliers	138981	3069 (2.70-2.70)
Sidechain outliers	138945	3069 (2.70-2.70)
RSRZ outliers	127900	2737 (2.70-2.70)

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	291	73% 19% • 6%
1	B	291	76% 17% • 6%
2	H	238	4% 82% 13% 5%
2	X	238	5% 80% 14% • 5%
3	L	216	% 78% 20% •

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Mol	Chain	Length	Quality of chain
3	Y	216	76% 22%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 11297 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 Lipoprotein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	273	2189	1388	365	429	7	0	0	0
1	B	274	2192	1388	366	431	7	0	0	0

- Molecule 2 is a protein called Heavy chain of antagonistic fab fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
2	H	227	1700	1074	279	341	6	0	0	0
2	X	227	1700	1074	279	341	6	0	0	0

- Molecule 3 is a protein called Light chain of antagonistic fab fragment.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
3	L	213	1634	1023	273	333	5	0	0	0
3	Y	213	1634	1023	273	333	5	0	0	0

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	63	63	63	0	0
4	B	49	49	49	0	0
4	H	27	27	27	0	0
4	X	35	35	35	0	0

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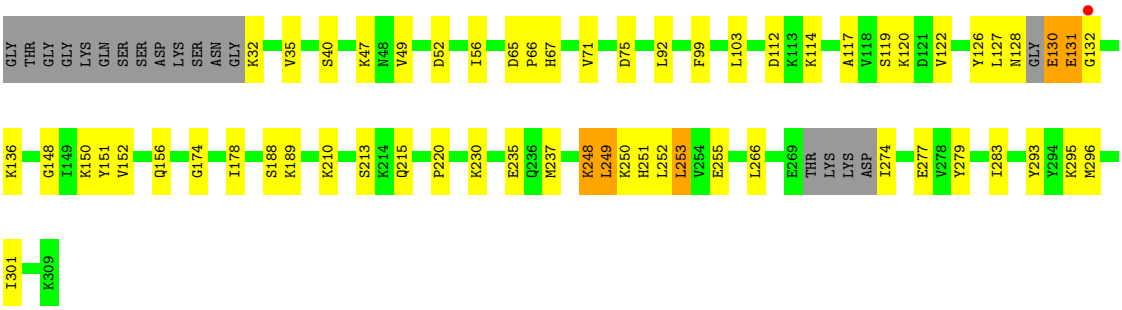
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	L	42	42	42	0	0
			Total	O		
4	Y	32	32	32	0	0

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.

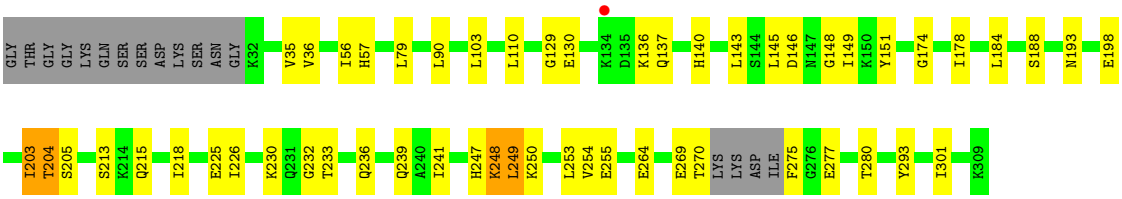
• Molecule 1: Lipoprotein

Chain A: 73% 19% • 6%



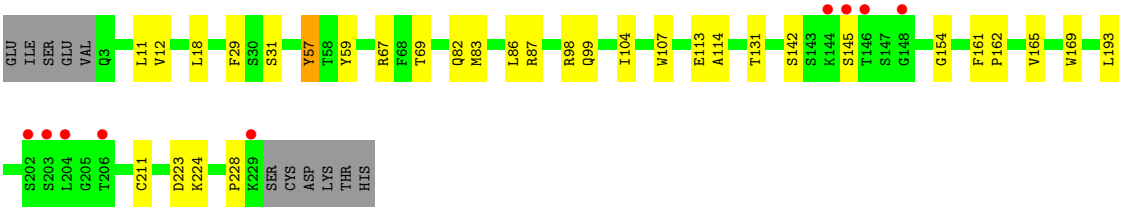
• Molecule 1: Lipoprotein

Chain B: 76% 17% • 6%



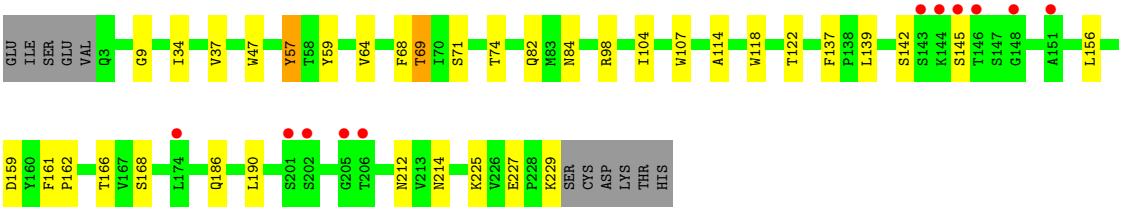
• Molecule 2: Heavy chain of antagonistic fab fragment

Chain H: 4% 82% 13% 5%



• Molecule 2: Heavy chain of antagonistic fab fragment

Chain X: 5% 80% 14% • 5%

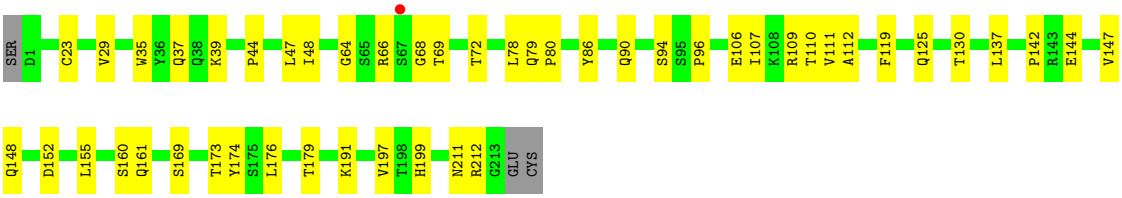


• Molecule 3: Light chain of antagonistic fab fragment

Chain L:

• Molecule 3: Light chain of antagonistic fab fragment

Chain Y:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	100.11Å 92.51Å 127.86Å 90.00° 92.48° 90.00°	Depositor
Resolution (Å)	47.27 – 2.69 47.27 – 2.68	Depositor EDS
% Data completeness (in resolution range)	88.9 (47.27-2.69) 88.9 (47.27-2.68)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.35 (at 2.69Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.224 , 0.283 0.225 , 0.281	Depositor DCC
R_{free} test set	2486 reflections (4.27%)	wwPDB-VP
Wilson B-factor (Å^2)	44.9	Xtriage
Anisotropy	0.605	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.33 , 37.6	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.027 for h,-k,-l	Xtriage
F_o,F_c correlation	0.91	EDS
Total number of atoms	11297	wwPDB-VP
Average B, all atoms (Å^2)	22.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 16.63% 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)

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	0.52	0/2227	0.72	2/2987 (0.1%)
1	B	0.53	0/2231	0.70	1/2994 (0.0%)
2	H	0.54	0/1744	0.65	0/2377
2	X	0.49	0/1744	0.64	0/2377
3	L	0.48	0/1670	0.63	0/2267
3	Y	0.50	0/1670	0.64	0/2267
All	All	0.51	0/11286	0.67	3/15269 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	249	LEU	CA-CB-CG	7.24	131.95	115.30
1	B	249	LEU	CA-CB-CG	6.14	129.42	115.30
1	A	252	LEU	CA-CB-CG	5.38	127.68	115.30

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	2189	0	2192	52	0
1	B	2192	0	2192	47	2
2	H	1700	0	1634	29	0
2	X	1700	0	1634	30	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	L	1634	0	1588	39	0
3	Y	1634	0	1588	37	0
4	A	63	0	0	2	0
4	B	49	0	0	1	0
4	H	27	0	0	2	2
4	L	42	0	0	1	0
4	X	35	0	0	1	0
4	Y	32	0	0	2	0
All	All	11297	0	10828	215	2

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 (215) 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:A:130:GLU:HA	1:A:130:GLU:OE1	1.48	1.08
1:B:255:GLU:HG3	1:B:277:GLU:OE1	1.54	1.06
1:B:253:LEU:O	1:B:253:LEU:HD12	1.62	1.00
1:A:266:LEU:HD11	3:L:27:GLN:HE22	1.31	0.92
1:A:128:ASN:C	1:A:130:GLU:HB2	1.93	0.88
1:B:255:GLU:HG3	1:B:277:GLU:CD	1.94	0.86
3:Y:109:ARG:HG2	3:Y:110:THR:H	1.42	0.85
2:H:87:ARG:NH2	4:H:302:HOH:O	2.09	0.85
1:B:253:LEU:HD12	1:B:253:LEU:C	2.03	0.79
1:A:99:PHE:O	1:A:103:LEU:HG	1.83	0.78
2:X:225:LYS:NZ	2:X:227:GLU:HB2	1.99	0.78
3:Y:68:GLY:O	3:Y:69:THR:HG22	1.84	0.76
2:H:224:LYS:NZ	3:L:124:GLU:HG2	2.00	0.76
1:B:204:THR:O	1:B:205:SER:CB	2.30	0.74
2:H:224:LYS:HZ1	3:L:124:GLU:HG2	1.53	0.74
2:X:168:SER:OG	2:X:212:ASN:HB2	1.87	0.73
2:H:69:THR:HG22	2:H:82:GLN:HB3	1.71	0.73
3:Y:109:ARG:HG2	3:Y:110:THR:N	2.04	0.72
2:X:166:THR:HG22	2:X:214:ASN:HB3	1.72	0.72
1:A:189:LYS:HE3	1:A:215:GLN:HE21	1.54	0.71
1:B:204:THR:O	1:B:205:SER:HB3	1.90	0.71
3:Y:191:LYS:HA	3:Y:212:ARG:HD3	1.72	0.70
1:B:193:ASN:ND2	4:B:422:HOH:O	2.23	0.70
1:B:253:LEU:HD11	1:B:277:GLU:HG2	1.72	0.70
2:X:166:THR:CG2	2:X:214:ASN:HB3	2.21	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:6:GLN:NE2	3:L:86:TYR:O	2.20	0.69
1:B:36:VAL:HG12	1:B:57:HIS:HB3	1.78	0.66
3:L:66:ARG:HG3	3:L:66:ARG:HH11	1.61	0.66
3:L:1:ASP:OD1	3:L:1:ASP:N	2.30	0.65
1:B:232:GLY:H	1:B:236:GLN:HE22	1.45	0.65
2:X:186:GLN:HA	3:Y:161:GLN:HE22	1.63	0.64
3:L:122:SER:OG	3:L:124:GLU:OE1	2.12	0.64
3:Y:152:ASP:HB3	4:Y:320:HOH:O	1.96	0.64
1:A:71:VAL:O	4:A:428:HOH:O	2.15	0.63
1:A:130:GLU:CG	1:A:136:LYS:CD	2.77	0.63
1:B:203:ILE:O	1:B:203:ILE:HG23	1.98	0.63
3:L:109:ARG:HG2	3:L:110:THR:H	1.64	0.62
2:H:69:THR:CG2	2:H:82:GLN:HB3	2.29	0.62
2:H:31:SER:OG	2:H:98:ARG:HD3	2.00	0.62
3:L:66:ARG:HG2	3:L:68:GLY:H	1.63	0.62
2:X:225:LYS:NZ	2:X:227:GLU:OE2	2.30	0.62
2:H:12:VAL:HG11	2:H:18:LEU:HD22	1.83	0.61
3:Y:109:ARG:CG	3:Y:110:THR:H	2.12	0.61
2:X:142:SER:H	2:X:145:SER:HB2	1.66	0.61
2:H:11:LEU:CD2	2:H:131:THR:HG23	2.31	0.61
1:B:36:VAL:HG21	1:B:79:LEU:HD23	1.83	0.60
3:L:35:TRP:HD1	3:L:48:ILE:HD11	1.67	0.60
3:Y:211:ASN:O	3:Y:212:ARG:HD2	2.02	0.59
1:B:230:LYS:HG3	2:X:107:TRP:CE3	2.37	0.59
3:Y:66:ARG:HG2	3:Y:68:GLY:H	1.68	0.59
1:A:130:GLU:HG2	1:A:136:LYS:HE3	1.85	0.59
1:B:140:HIS:HB3	1:B:143:LEU:HD12	1.85	0.58
1:A:130:GLU:HG3	1:A:136:LYS:HD2	1.85	0.57
3:L:19:VAL:HG21	3:L:78:LEU:HD22	1.87	0.57
1:B:203:ILE:O	1:B:203:ILE:CG2	2.54	0.56
3:L:66:ARG:HG3	3:L:66:ARG:NH1	2.21	0.56
2:H:104:ILE:HD12	2:H:114:ALA:HB3	1.88	0.55
2:X:225:LYS:HZ2	2:X:227:GLU:HB2	1.69	0.55
2:X:142:SER:N	2:X:145:SER:HB2	2.21	0.55
1:B:233:THR:H	1:B:236:GLN:NE2	2.04	0.55
3:Y:109:ARG:CG	3:Y:110:THR:N	2.70	0.54
1:B:188:SER:HB3	1:B:301:ILE:HD13	1.89	0.54
3:Y:137:LEU:HD13	3:Y:176:LEU:HD22	1.88	0.54
3:Y:106:GLU:HG3	3:Y:174:TYR:OH	2.08	0.54
2:H:142:SER:H	2:H:145:SER:HB2	1.72	0.54
2:X:159:ASP:HB3	2:X:190:LEU:HD13	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:11:LEU:HD22	2:H:131:THR:CG2	2.38	0.54
2:X:64:VAL:HB	2:X:68:PHE:CG	2.42	0.54
1:B:148:GLY:O	1:B:151:TYR:HB2	2.08	0.54
2:X:104:ILE:HD12	2:X:114:ALA:HB3	1.90	0.54
3:Y:66:ARG:CG	3:Y:68:GLY:H	2.21	0.54
3:Y:109:ARG:HH21	3:Y:173:THR:HG22	1.73	0.54
3:L:109:ARG:HG2	3:L:110:THR:N	2.22	0.54
1:B:35:VAL:HG12	1:B:56:ILE:HG23	1.91	0.53
1:A:148:GLY:O	1:A:151:TYR:HB2	2.07	0.53
1:A:130:GLU:HG3	1:A:136:LYS:CD	2.39	0.53
1:A:295:LYS:NZ	4:A:462:HOH:O	2.42	0.52
1:B:236:GLN:O	1:B:239:GLN:HG3	2.10	0.52
1:B:255:GLU:CG	1:B:277:GLU:CD	2.72	0.52
3:L:66:ARG:CG	3:L:68:GLY:H	2.22	0.52
3:L:106:GLU:HG2	3:L:107:ILE:N	2.25	0.51
1:B:232:GLY:H	1:B:236:GLN:NE2	2.07	0.51
2:H:228:PRO:HD3	4:H:313:HOH:O	2.09	0.51
2:H:99:GLN:HE22	2:H:113:GLU:HB3	1.75	0.51
2:X:57:TYR:CD2	2:X:59:TYR:CZ	2.99	0.51
1:A:251:HIS:HB2	1:A:274:ILE:HA	1.91	0.50
3:Y:68:GLY:O	3:Y:69:THR:CG2	2.55	0.50
3:L:37:GLN:CD	3:L:47:LEU:HD21	2.31	0.50
3:Y:78:LEU:HD12	3:Y:79:GLN:H	1.77	0.50
1:A:67:HIS:HA	1:A:92:LEU:O	2.12	0.50
1:A:130:GLU:CG	1:A:136:LYS:HD3	2.41	0.50
3:Y:125:GLN:HG2	3:Y:130:THR:O	2.12	0.49
1:A:119:SER:O	1:A:122:VAL:HG13	2.13	0.49
3:L:148:GLN:HG2	3:L:155:LEU:HD11	1.95	0.49
2:H:131:THR:HG22	2:H:162:PRO:HD3	1.94	0.49
1:B:264:GLU:HG2	1:B:269:GLU:O	2.12	0.49
3:L:2:ILE:HG12	3:L:27:GLN:HG2	1.94	0.49
3:Y:94:SER:N	4:Y:319:HOH:O	2.27	0.49
2:H:211:CYS:O	2:H:223:ASP:HA	2.13	0.48
1:A:132:GLY:HA3	2:X:71:SER:HB2	1.95	0.48
3:L:35:TRP:HB2	3:L:48:ILE:HG13	1.94	0.48
1:B:248:LYS:O	1:B:249:LEU:HD23	2.14	0.48
3:L:109:ARG:NH1	3:L:112:ALA:HB2	2.29	0.48
3:L:37:GLN:NE2	4:L:321:HOH:O	2.46	0.48
1:A:255:GLU:HB2	1:A:277:GLU:HB3	1.96	0.48
3:L:188:GLU:O	3:L:212:ARG:NH1	2.47	0.48
3:Y:144:GLU:O	3:Y:199:HIS:HD2	1.96	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:188:SER:CB	1:B:301:ILE:HD13	2.43	0.48
2:H:57:TYR:CD2	2:H:59:TYR:CZ	3.02	0.48
1:B:205:SER:HB3	1:B:254:VAL:HG23	1.96	0.47
1:B:253:LEU:O	1:B:253:LEU:CD1	2.48	0.47
1:B:145:LEU:HD11	1:B:184:LEU:HD23	1.96	0.47
3:L:156:GLN:HG3	3:L:180:LEU:HD11	1.95	0.47
1:A:150:LYS:HE2	1:B:129:GLY:O	2.14	0.47
1:B:145:LEU:O	1:B:149:ILE:HG13	2.14	0.47
1:A:131:GLU:HG3	2:X:74:THR:HG23	1.96	0.47
1:A:213:SER:OG	1:A:220:PRO:HB3	2.14	0.47
2:X:225:LYS:HZ3	2:X:227:GLU:HB2	1.77	0.47
3:Y:68:GLY:C	3:Y:69:THR:HG22	2.36	0.46
1:B:146:ASP:OD2	1:B:215:GLN:HG3	2.15	0.46
2:H:99:GLN:NE2	2:H:113:GLU:HB3	2.31	0.46
3:L:211:ASN:O	3:L:213:GLY:N	2.46	0.46
3:Y:47:LEU:O	3:Y:48:ILE:HD12	2.15	0.46
3:Y:80:PRO:HB2	3:Y:169:SER:O	2.15	0.46
1:B:270:THR:O	1:B:275:PHE:N	2.48	0.46
3:L:64:GLY:HA2	3:L:72:THR:O	2.16	0.46
1:A:249:LEU:HB3	1:A:250:LYS:H	1.48	0.46
1:B:225:GLU:OE2	1:B:226:ILE:HG12	2.15	0.46
3:Y:106:GLU:HG2	3:Y:107:ILE:N	2.31	0.46
3:L:202:LEU:HD13	3:L:206:VAL:HG23	1.98	0.45
3:Y:37:GLN:HB2	3:Y:86:TYR:CE2	2.52	0.45
1:B:174:GLY:O	1:B:178:ILE:HG13	2.17	0.45
1:B:148:GLY:HA3	1:B:293:TYR:OH	2.16	0.45
2:H:12:VAL:CG1	2:H:18:LEU:HD13	2.47	0.45
1:B:247:HIS:O	1:B:248:LYS:HB3	2.16	0.45
2:X:69:THR:HG22	4:X:310:HOH:O	2.17	0.45
3:L:18:ARG:HG3	3:L:76:SER:HA	1.99	0.45
2:H:224:LYS:HZ3	3:L:124:GLU:HG2	1.78	0.45
1:B:130:GLU:OE1	1:B:136:LYS:HG3	2.16	0.45
3:L:200:GLN:O	3:L:200:GLN:HG2	2.17	0.45
1:A:127:LEU:CD2	1:A:210:LYS:NZ	2.80	0.44
2:H:83:MET:HE2	2:H:86:LEU:HD21	1.99	0.44
3:Y:39:LYS:HB2	3:Y:39:LYS:HE3	1.71	0.44
3:Y:64:GLY:HA2	3:Y:72:THR:O	2.17	0.44
1:A:99:PHE:CE2	1:A:103:LEU:HD11	2.51	0.44
1:A:230:LYS:HG3	2:H:107:TRP:CZ3	2.52	0.44
1:B:241:ILE:HD13	3:Y:94:SER:HA	1.99	0.44
3:Y:148:GLN:HG2	3:Y:155:LEU:HD11	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:112:ASP:HB3	1:A:114:LYS:NZ	2.32	0.44
1:A:130:GLU:HG2	1:A:136:LYS:CE	2.48	0.43
1:A:266:LEU:HD11	3:L:27:GLN:NE2	2.14	0.43
1:A:47:LYS:HD2	1:A:47:LYS:HA	1.80	0.43
1:A:174:GLY:O	1:A:178:ILE:HG13	2.18	0.43
1:B:253:LEU:CD1	1:B:255:GLU:HB2	2.48	0.43
3:Y:109:ARG:HH22	3:Y:112:ALA:HB2	1.84	0.43
1:A:65:ASP:HA	1:A:66:PRO:HD3	1.92	0.43
1:A:248:LYS:O	1:A:249:LEU:HD23	2.18	0.43
1:A:279:TYR:HB2	1:A:296:MET:HG3	2.01	0.43
3:Y:147:VAL:HG22	3:Y:197:VAL:HG22	2.00	0.43
2:X:82:GLN:NE2	2:X:84:ASN:OD1	2.46	0.43
2:H:29:PHE:CZ	2:H:98:ARG:HD2	2.53	0.43
2:X:37:VAL:HG22	2:X:47:TRP:HA	2.01	0.43
3:L:66:ARG:HB3	3:L:70:ASP:O	2.19	0.43
1:A:131:GLU:HB2	2:X:74:THR:N	2.34	0.42
1:B:230:LYS:HG3	2:X:107:TRP:CZ3	2.54	0.42
1:B:249:LEU:HB3	1:B:250:LYS:H	1.48	0.42
2:H:67:ARG:HH11	2:H:67:ARG:HD2	1.72	0.42
1:A:99:PHE:O	1:A:103:LEU:CG	2.62	0.42
1:A:130:GLU:HG2	1:A:136:LYS:CD	2.48	0.42
1:A:71:VAL:HG13	1:A:75:ASP:HB2	2.00	0.42
1:B:103:LEU:HD23	1:B:103:LEU:HA	1.91	0.42
2:H:165:VAL:CG1	2:H:193:LEU:HD21	2.49	0.42
1:A:32:LYS:HE3	1:A:52:ASP:O	2.20	0.42
1:A:117:ALA:O	1:A:120:LYS:HB3	2.19	0.42
1:B:213:SER:HB3	1:B:218:ILE:O	2.20	0.42
2:X:137:PHE:HB2	2:X:156:LEU:HB3	2.02	0.42
3:L:78:LEU:HD12	3:L:79:GLN:H	1.84	0.42
3:Y:111:VAL:HG22	3:Y:142:PRO:HD3	2.01	0.42
2:H:12:VAL:CG1	2:H:18:LEU:HD22	2.48	0.42
1:A:148:GLY:HA3	1:A:293:TYR:OH	2.19	0.42
2:X:139:LEU:HB3	3:Y:119:PHE:CD1	2.55	0.42
1:B:143:LEU:HD21	1:B:280:THR:HA	2.02	0.42
3:L:55:TYR:O	3:L:58:VAL:HB	2.20	0.42
1:A:130:GLU:CB	1:A:136:LYS:HD3	2.50	0.41
2:X:34:ILE:HG12	2:X:98:ARG:HG2	2.01	0.41
2:H:161:PHE:HA	2:H:162:PRO:HA	1.74	0.41
3:L:19:VAL:CG2	3:L:78:LEU:HD22	2.49	0.41
2:X:161:PHE:HA	2:X:162:PRO:HA	1.85	0.41
3:L:137:LEU:HD13	3:L:176:LEU:HD22	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:49:VAL:HG11	1:A:156:GLN:HA	2.03	0.41
1:A:230:LYS:HG3	2:H:107:TRP:CE3	2.56	0.41
1:A:253:LEU:CD1	1:A:277:GLU:HG2	2.51	0.41
3:Y:160:SER:HA	3:Y:179:THR:O	2.20	0.41
1:B:90:LEU:HA	1:B:137:GLN:NE2	2.35	0.41
3:Y:29:VAL:HG11	3:Y:90:GLN:HG3	2.03	0.41
1:A:188:SER:HB3	1:A:301:ILE:HD13	2.03	0.41
1:B:204:THR:O	1:B:254:VAL:CG2	2.69	0.41
3:L:37:GLN:HB3	3:L:86:TYR:CE2	2.56	0.41
1:A:237:MET:HB3	3:L:93:TYR:CD1	2.55	0.41
2:H:154:GLY:HA2	2:H:169:TRP:CH2	2.56	0.41
2:X:118:TRP:CE3	3:Y:44:PRO:HD2	2.55	0.41
1:A:49:VAL:CG2	1:A:152:VAL:HG13	2.51	0.41
1:B:110:LEU:HD12	1:B:110:LEU:HA	1.89	0.41
2:H:11:LEU:CD2	2:H:131:THR:CG2	2.95	0.41
2:X:47:TRP:CZ3	3:Y:96:PRO:HB3	2.56	0.40
2:X:229:LYS:HD3	2:X:229:LYS:HA	1.81	0.40
3:L:35:TRP:CD2	3:L:73:LEU:HB2	2.56	0.40
3:Y:23:CYS:HB2	3:Y:35:TRP:CH2	2.56	0.40
1:A:40:SER:HB2	1:A:283:ILE:HD13	2.02	0.40
1:A:127:LEU:O	1:A:128:ASN:C	2.59	0.40
1:A:35:VAL:HG12	1:A:56:ILE:HG23	2.04	0.40
1:A:99:PHE:CZ	1:A:103:LEU:HD21	2.56	0.40
1:A:126:TYR:CG	1:A:131:GLU:HA	2.56	0.40
2:X:9:GLY:HA3	2:X:122:THR:OG1	2.21	0.40
3:L:125:GLN:HG2	3:L:130:THR:O	2.21	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:198:GLU:OE1	4:H:302:HOH:O[2_545]	1.21	0.99
1:B:198:GLU:CD	4:H:302:HOH:O[2_545]	2.03	0.17

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	267/291 (92%)	247 (92%)	17 (6%)	3 (1%)	14	34
1	B	270/291 (93%)	251 (93%)	18 (7%)	1 (0%)	34	60
2	H	225/238 (94%)	215 (96%)	10 (4%)	0	100	100
2	X	225/238 (94%)	214 (95%)	11 (5%)	0	100	100
3	L	211/216 (98%)	197 (93%)	13 (6%)	1 (0%)	29	54
3	Y	211/216 (98%)	193 (92%)	18 (8%)	0	100	100
All	All	1409/1490 (95%)	1317 (94%)	87 (6%)	5 (0%)	34	60

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	131	GLU
1	A	235	GLU
1	A	248	LYS
1	B	248	LYS
3	L	212	ARG

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	242/255 (95%)	240 (99%)	2 (1%)	81	93
1	B	242/255 (95%)	240 (99%)	2 (1%)	81	93
2	H	188/199 (94%)	187 (100%)	1 (0%)	88	96
2	X	188/199 (94%)	186 (99%)	2 (1%)	73	90
3	L	188/191 (98%)	188 (100%)	0	100	100
3	Y	188/191 (98%)	188 (100%)	0	100	100
All	All	1236/1290 (96%)	1229 (99%)	7 (1%)	86	95

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	130	GLU
1	A	253	LEU
1	B	203	ILE
1	B	204	THR
2	H	57	TYR
2	X	57	TYR
2	X	69	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	53	ASN
1	A	173	GLN
1	A	185	ASN
1	A	186	ASN
1	A	215	GLN
1	A	251	HIS
1	B	186	ASN
1	B	236	GLN
2	H	28	ASN
2	H	99	GLN
3	L	27	GLN
3	L	89	GLN
3	L	138	ASN
3	L	148	GLN
3	Y	27	GLN
3	Y	161	GLN

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)

There are no ligands in this entry.

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	273/291 (93%)	-0.22	1 (0%)	92	93	3, 14, 48, 96	0
1	B	274/291 (94%)	-0.14	1 (0%)	92	93	4, 15, 49, 69	0
2	H	227/238 (95%)	0.01	9 (3%)	38	37	3, 20, 78, 110	0
2	X	227/238 (95%)	0.04	11 (4%)	30	28	2, 19, 77, 118	0
3	L	213/216 (98%)	-0.17	2 (0%)	84	85	5, 20, 53, 64	0
3	Y	213/216 (98%)	-0.17	1 (0%)	91	92	3, 23, 55, 67	0
All	All	1427/1490 (95%)	-0.11	25 (1%)	68	70	2, 18, 57, 118	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	X	145	SER	5.6
2	X	148	GLY	4.8
2	H	146	THR	4.5
2	H	206	THR	3.8
2	H	145	SER	3.7
2	X	206	THR	3.5
2	X	146	THR	3.4
2	H	148	GLY	3.4
2	H	229	LYS	3.3
2	H	144	LYS	3.2
2	X	144	LYS	3.1
2	X	143	SER	3.1
1	A	132	GLY	2.9
2	X	151	ALA	2.7
2	H	204	LEU	2.5
1	B	134	LYS	2.5
2	H	202	SER	2.5
3	Y	67	SER	2.4
2	H	203	SER	2.4

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Mol	Chain	Res	Type	RSRZ
2	X	174	LEU	2.4
2	X	202	SER	2.2
2	X	205	GLY	2.2
3	L	37	GLN	2.1
2	X	201	SER	2.1
3	L	67	SER	2.1

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.