Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 09:05 AM GMT

PDB ID	:	3H42
Title	:	Crystal structure of PCSK9 in complex with Fab from LDLR competitive antibody
Authors	:	Piper, D.E.; Walker, N.P.C.; Romanow, W.G.; Thibault, S.T.; Tsai, M.M.; Yang, E.
Deposited on	:	2009-04-17
Resolution	:	2.30 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
http://wwpdb.org/validation/2016/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The following versions of software and data (see references (1)) were used in the production of this report:

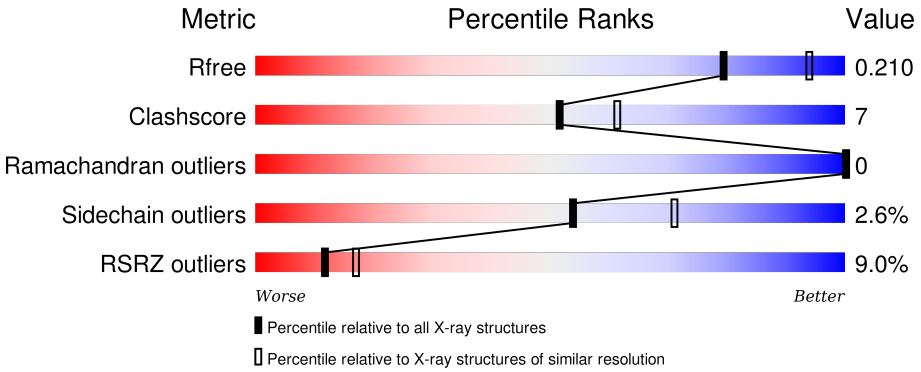
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	trunk26865

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.30 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}	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (2.30-2.30)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments on the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	126	3%, 58%, 14%, 27%
2	B	540	13%, 76%, 14%, 9%
3	L	217	5%, 87%, 12%
4	H	238	3%, 82%, 11%, 7%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
5	NA	B	1	-	-	-	X

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 8196 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 Proprotein convertase subtilisin/kexin type 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	92	Total	C	N	O	S	0	0	0
			740	474	133	131	2			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	27	GLY	-	EXPRESSION TAG	UNP Q8NBP7
A	28	ALA	-	EXPRESSION TAG	UNP Q8NBP7
A	29	MET	-	EXPRESSION TAG	UNP Q8NBP7
A	30	GLY	-	EXPRESSION TAG	UNP Q8NBP7

- Molecule 2 is a protein called Proprotein convertase subtilisin/kexin type 9.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	492	Total	C	N	O	S	0	0	0
			3659	2261	676	690	32			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	533	ALA	ASN	ENGINEERED	UNP Q8NBP7
B	620	GLY	GLU	SEE REMARK 999	UNP Q8NBP7

- Molecule 3 is a protein called Fab from LDLR competitive antibody: Light chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	L	214	Total	C	N	O	S	0	0	0
			1568	975	263	326	4			

- Molecule 4 is a protein called Fab from LDLR competitive antibody: Heavy chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	H	221	Total	C	N	O	S	0	0	0
			1660	1051	274	328	7			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
H	233	HIS	-	EXPRESSION TAG	PDB 3H42
H	234	HIS	-	EXPRESSION TAG	PDB 3H42
H	235	HIS	-	EXPRESSION TAG	PDB 3H42
H	236	HIS	-	EXPRESSION TAG	PDB 3H42
H	237	HIS	-	EXPRESSION TAG	PDB 3H42
H	238	HIS	-	EXPRESSION TAG	PDB 3H42

- Molecule 5 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	1	Total	Na	0	0
			1	1		

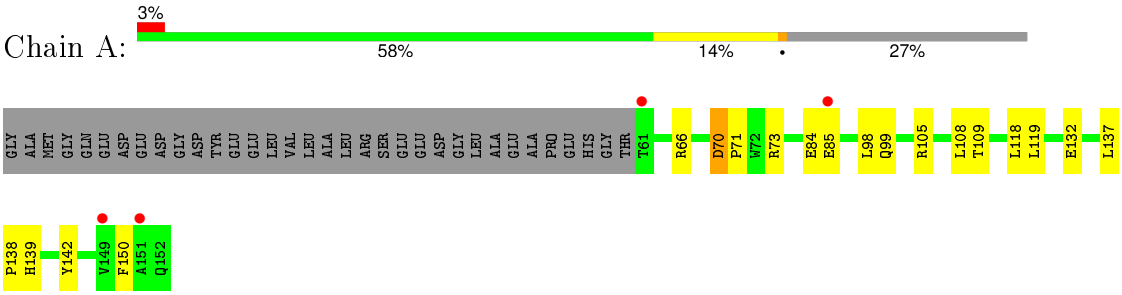
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	76	Total	O	0	0
			76	76		
6	B	178	Total	O	0	0
			178	178		
6	L	159	Total	O	0	0
			159	159		
6	H	155	Total	O	0	0
			155	155		

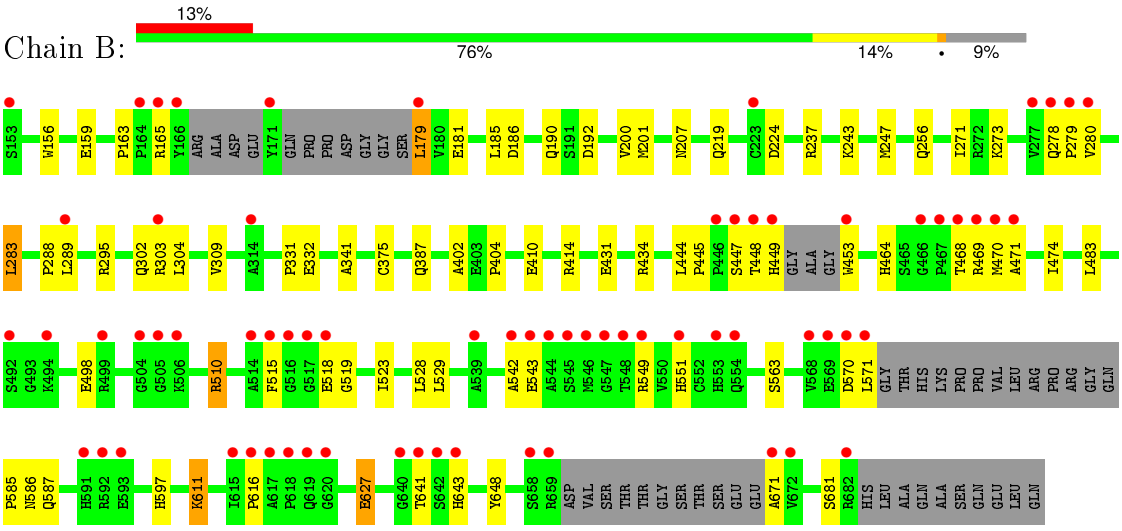
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

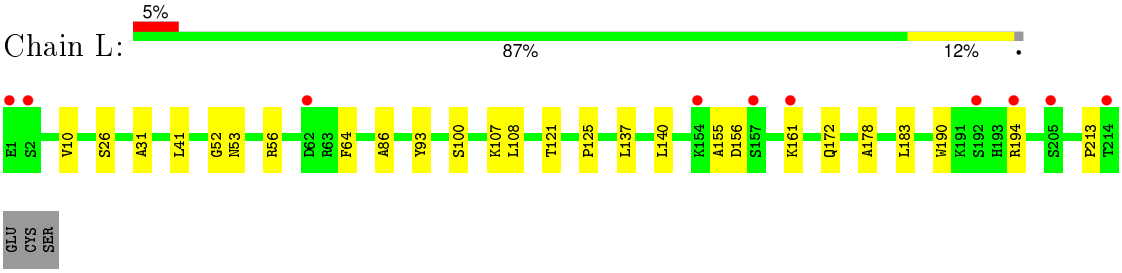
- Molecule 1: Proprotein convertase subtilisin/kexin type 9



- Molecule 2: Proprotein convertase subtilisin/kexin type 9

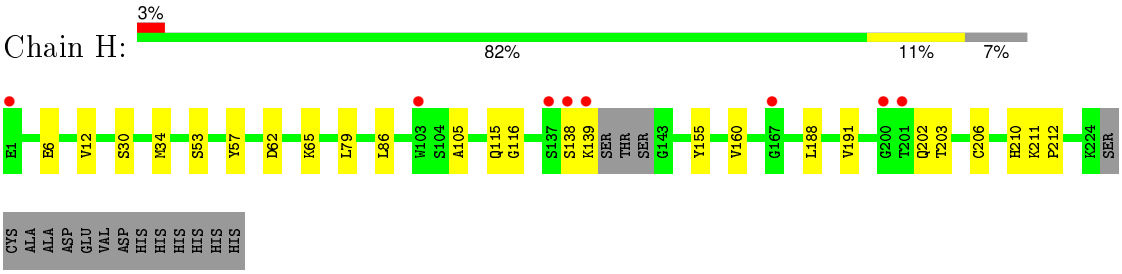


- Molecule 3: Fab from LDLR competitive antibody: Light chain



- Molecule 4: Fab from LDLR competitive antibody: Heavy chain

Chain H:


3%
82%
11%
7%
E1
E6
V12
S30
M34
S53
Y57
D62
K65
L79
L86
W103
S104
A105
Q115
G116
S137
S138
K139
SER
THR
SER
G143
Y155
V160
G167
L188
V191
G200
T201
Q202
T203
C206
H210
K211
P212
K224
SER
CYS
ALA
ALA
ASP
GLU
VAL
ASP
HIS
HIS
HIS
HIS
HIS
HIS

4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α, β, γ	264.72Å 137.35Å 69.89Å 90.00° 102.84° 90.00°	Depositor
Resolution (Å)	40.00 – 2.30 38.29 – 2.30	Depositor EDS
% Data completeness (in resolution range)	99.9 (40.00-2.30) 99.8 (38.29-2.30)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.53 (at 2.29Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.191 , 0.209 0.191 , 0.210	Depositor DCC
R_{free} test set	5376 reflections (4.99%)	DCC
Wilson B-factor (Å^2)	38.0	Xtriage
Anisotropy	0.481	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.35 , 49.5	EDS
Estimated twinning fraction	0.022 for -h-2*l,-k,l	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.32$	Xtriage
Outliers	0 of 107765 reflections	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	8196	wwPDB-VP
Average B, all atoms (Å^2)	46.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.50% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

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

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	0.31	0/757	0.43	0/1023
2	B	0.26	0/3729	0.47	0/5064
3	L	0.30	0/1605	0.49	0/2192
4	H	0.32	0/1701	0.49	0/2313
All	All	0.29	0/7792	0.47	0/10592

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

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	740	0	750	15	0
2	B	3659	0	3572	56	0
3	L	1568	0	1518	18	0
4	H	1660	0	1610	18	0
5	B	1	0	0	0	0
6	A	76	0	0	2	0
6	B	178	0	0	1	0
6	H	155	0	0	1	0
6	L	159	0	0	2	0
All	All	8196	0	7450	103	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 7.

All (103) 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:B:474:ILE:HD11	2:B:510:ARG:HH21	1.40	0.84
2:B:200:VAL:HG12	2:B:247:MET:HB2	1.60	0.82
1:A:66:ARG:HH21	1:A:73:ARG:HD2	1.43	0.81
2:B:474:ILE:HD11	2:B:510:ARG:NH2	1.97	0.80
2:B:402:ALA:HB1	2:B:448:THR:HG23	1.65	0.77
2:B:414:ARG:HH11	2:B:414:ARG:HG3	1.50	0.74
2:B:468:THR:HB	2:B:471:ALA:HB2	1.71	0.73
2:B:469:ARG:HB2	2:B:515:PHE:HD2	1.58	0.68
2:B:278:GLN:HB3	2:B:279:PRO:HD2	1.76	0.68
2:B:302:GLN:HA	2:B:332:GLU:HG3	1.76	0.67
2:B:185:LEU:HD11	2:B:271:ILE:HD11	1.78	0.66
2:B:186:ASP:OD1	2:B:288:PRO:HG2	1.97	0.65
2:B:332:GLU:CD	2:B:332:GLU:H	2.00	0.65
2:B:523:ILE:HD13	2:B:648:TYR:HB3	1.80	0.63
3:L:125:PRO:HD3	3:L:137:LEU:CD2	2.28	0.63
1:A:84:GLU:O	1:A:85:GLU:HB2	2.01	0.61
4:H:138:SER:C	4:H:139:LYS:HD2	2.21	0.60
2:B:570:ASP:C	2:B:571:LEU:HD12	2.23	0.59
1:A:66:ARG:NH2	1:A:73:ARG:HD2	2.17	0.58
2:B:304:LEU:HD22	2:B:309:VAL:HG21	1.86	0.58
2:B:641:THR:HG22	2:B:643:HIS:H	1.69	0.57
2:B:414:ARG:NH1	2:B:414:ARG:HG3	2.18	0.56
2:B:468:THR:HG22	2:B:470:MET:H	1.71	0.56
2:B:448:THR:O	2:B:449:HIS:HB2	2.05	0.55
2:B:179:LEU:HD13	2:B:404:PRO:HG3	1.89	0.54
2:B:611:LYS:HB3	2:B:611:LYS:NZ	2.23	0.54
3:L:108:LEU:HD23	3:L:108:LEU:C	2.28	0.54
4:H:34:MET:HB3	4:H:79:LEU:HD22	1.91	0.53
4:H:160:VAL:CG2	4:H:188:LEU:HD21	2.38	0.53
2:B:256:GLN:HB3	4:H:57:TYR:CZ	2.44	0.53
3:L:125:PRO:HD3	3:L:137:LEU:HD21	1.91	0.52
2:B:474:ILE:CD1	2:B:510:ARG:HH21	2.20	0.51
2:B:331:PRO:HD2	2:B:332:GLU:OE2	2.10	0.51
4:H:202:GLN:HE21	4:H:203:THR:N	2.08	0.51
2:B:445:PRO:C	2:B:447:SER:H	2.13	0.51
1:A:118:LEU:O	1:A:119:LEU:HD23	2.12	0.50
3:L:140:LEU:CD1	4:H:191:VAL:HG21	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:142:TYR:CE1	2:B:295:ARG:HG2	2.47	0.49
4:H:160:VAL:HG22	4:H:188:LEU:HD21	1.95	0.49
1:A:132:GLU:HG2	6:A:177:HOH:O	2.13	0.48
2:B:283:LEU:HD21	2:B:309:VAL:HG22	1.95	0.48
2:B:464:HIS:NE2	2:B:519:GLY:HA3	2.28	0.48
2:B:563:SER:HB2	2:B:597:HIS:HB2	1.96	0.48
4:H:160:VAL:HG12	4:H:210:HIS:CD2	2.48	0.48
2:B:190:GLN:NE2	2:B:192:ASP:HB2	2.28	0.48
4:H:6:GLU:CD	4:H:116:GLY:H	2.17	0.48
2:B:542:ALA:O	2:B:543:GLU:HB2	2.14	0.47
3:L:155:ALA:O	3:L:156:ASP:HB2	2.15	0.47
3:L:26:SER:O	3:L:31:ALA:HB2	2.15	0.47
2:B:283:LEU:H	2:B:283:LEU:HD23	1.79	0.47
3:L:56:ARG:HD3	3:L:64:PHE:O	2.14	0.47
3:L:156:ASP:OD1	3:L:194:ARG:HB2	2.16	0.46
3:L:137:LEU:HD12	3:L:183:LEU:HD23	1.98	0.46
2:B:224:ASP:HB2	6:B:759:HOH:O	2.16	0.46
3:L:41:LEU:HD23	3:L:86:ALA:HB2	1.98	0.45
2:B:156:TRP:CZ3	2:B:341:ALA:HA	2.50	0.45
4:H:115:GLN:NE2	6:H:461:HOH:O	2.49	0.45
4:H:188:LEU:HD12	4:H:188:LEU:C	2.37	0.45
2:B:190:GLN:HE21	2:B:192:ASP:H	1.64	0.45
2:B:387:GLN:NE2	2:B:387:GLN:H	2.15	0.45
3:L:125:PRO:HD3	3:L:137:LEU:HD23	1.98	0.44
2:B:611:LYS:HD3	2:B:627:GLU:OE2	2.18	0.44
4:H:202:GLN:HE21	4:H:203:THR:H	1.64	0.44
4:H:211:LYS:N	4:H:212:PRO:CD	2.81	0.44
1:A:105:ARG:HG3	1:A:105:ARG:HH21	1.82	0.44
1:A:84:GLU:O	1:A:85:GLU:CB	2.65	0.44
1:A:150:PHE:CG	4:H:105:ALA:HB2	2.53	0.44
2:B:237:ARG:O	2:B:243:LYS:HD3	2.19	0.43
2:B:571:LEU:HD12	2:B:571:LEU:N	2.32	0.43
1:A:138:PRO:O	1:A:139:HIS:HB2	2.19	0.43
2:B:453:TRP:CZ3	2:B:681:SER:HB2	2.53	0.43
1:A:118:LEU:HD12	1:A:118:LEU:HA	1.91	0.43
3:L:190:TRP:CZ2	3:L:213:PRO:HA	2.54	0.43
1:A:70:ASP:N	1:A:71:PRO:HD2	2.33	0.43
4:H:12:VAL:HG11	4:H:86:LEU:HD13	2.00	0.43
4:H:65:LYS:HB2	4:H:65:LYS:HE2	1.77	0.42
2:B:616:PRO:HA	2:B:671:ALA:HB2	2.01	0.42
2:B:431:GLU:HA	2:B:434:ARG:HD2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:10:VAL:HG12	6:L:456:HOH:O	2.20	0.42
3:L:93:TYR:HA	3:L:100:SER:HA	2.00	0.42
2:B:469:ARG:HB2	2:B:515:PHE:CD2	2.46	0.42
2:B:273:LYS:HE2	2:B:273:LYS:HB2	1.86	0.42
3:L:161:LYS:HB2	6:L:484:HOH:O	2.19	0.41
1:A:108:LEU:HD12	6:A:167:HOH:O	2.20	0.41
2:B:185:LEU:O	2:B:288:PRO:HD2	2.21	0.41
3:L:172:GLN:OE1	3:L:178:ALA:HB2	2.21	0.41
2:B:289:LEU:HD12	2:B:289:LEU:C	2.40	0.41
2:B:549:ARG:HH11	2:B:587:GLN:HE22	1.67	0.41
2:B:414:ARG:NH1	2:B:414:ARG:CG	2.82	0.41
1:A:98:LEU:HB2	1:A:137:LEU:HD11	2.02	0.41
2:B:302:GLN:HA	2:B:332:GLU:CG	2.47	0.41
3:L:194:ARG:HA	3:L:194:ARG:HD3	1.92	0.41
4:H:30:SER:O	4:H:53:SER:HB2	2.21	0.41
2:B:185:LEU:N	2:B:185:LEU:HD12	2.36	0.41
2:B:585:PRO:O	2:B:586:ASN:HB2	2.20	0.41
2:B:280:VAL:O	2:B:280:VAL:HG23	2.21	0.41
2:B:163:PRO:HG3	2:B:444:LEU:O	2.22	0.40
2:B:469:ARG:HG3	2:B:470:MET:HG2	2.04	0.40
1:A:99:GLN:HG2	1:A:109:THR:OG1	2.21	0.40
3:L:52:GLY:O	3:L:53:ASN:HB2	2.22	0.40
4:H:155:TYR:CE1	4:H:160:VAL:HG13	2.57	0.40
2:B:551:HIS:HB2	2:B:586:ASN:O	2.22	0.40
2:B:410:GLU:HA	2:B:528:LEU:HD11	2.04	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	90/126 (71%)	87 (97%)	3 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	481/540 (89%)	466 (97%)	15 (3%)	0	100	100
3	L	212/217 (98%)	207 (98%)	5 (2%)	0	100	100
4	H	217/238 (91%)	216 (100%)	1 (0%)	0	100	100
All	All	1000/1121 (89%)	976 (98%)	24 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	79/104 (76%)	78 (99%)	1 (1%)	76	87
2	B	392/429 (91%)	375 (96%)	17 (4%)	35	47
3	L	177/180 (98%)	175 (99%)	2 (1%)	80	90
4	H	186/201 (92%)	184 (99%)	2 (1%)	80	90
All	All	834/914 (91%)	812 (97%)	22 (3%)	54	71

All (22) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	70	ASP
2	B	159	GLU
2	B	165	ARG
2	B	179	LEU
2	B	181	GLU
2	B	201	MET
2	B	207	ASN
2	B	219	GLN
2	B	283	LEU
2	B	303	ARG
2	B	375	CYS
2	B	483	LEU
2	B	498	GLU

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Mol	Chain	Res	Type
2	B	510	ARG
2	B	518	GLU
2	B	529	LEU
2	B	611	LYS
2	B	627	GLU
3	L	107	LYS
3	L	121	THR
4	H	62	ASP
4	H	206	CYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	101	GLN
2	B	190	GLN
2	B	207	ASN
2	B	217	HIS
2	B	382	GLN
2	B	387	GLN
2	B	454	GLN
2	B	553	HIS
2	B	587	GLN
4	H	3	GLN
4	H	202	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	92/126 (73%)	0.13	4 (4%) 39 48	22, 39, 56, 69	0
2	B	492/540 (91%)	0.87	70 (14%) 4 6	23, 48, 93, 116	0
3	L	214/217 (98%)	-0.01	10 (4%) 35 44	23, 36, 61, 75	0
4	H	221/238 (92%)	0.16	8 (3%) 46 55	23, 38, 57, 91	0
All	All	1019/1121 (90%)	0.46	92 (9%) 12 17	22, 41, 83, 116	0

All (92) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	546	MET	10.3
2	B	515	PHE	8.6
2	B	453	TRP	7.5
2	B	545	SER	7.4
2	B	643	HIS	6.8
4	H	138	SER	6.4
2	B	469	ARG	6.4
2	B	165	ARG	6.4
2	B	171	TYR	6.0
2	B	277	VAL	6.0
2	B	619	GLN	5.8
2	B	642	SER	5.6
2	B	447	SER	5.4
2	B	166	TYR	5.4
2	B	153	SER	5.3
2	B	659	ARG	5.2
2	B	470	MET	5.2
2	B	616	PRO	5.0
3	L	194	ARG	4.7
3	L	214	THR	4.7
2	B	553	HIS	4.6

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Mol	Chain	Res	Type	RSRZ
2	B	641	THR	4.6
2	B	570	ASP	4.5
2	B	448	THR	4.5
2	B	471	ALA	4.4
2	B	164	PRO	4.4
2	B	446	PRO	4.3
2	B	615	ILE	4.3
2	B	617	ALA	4.3
2	B	618	PRO	4.2
2	B	547	GLY	4.1
1	A	85	GLU	4.0
2	B	682	ARG	4.0
2	B	449	HIS	3.9
2	B	640	GLY	3.9
2	B	592	ARG	3.8
2	B	544	ALA	3.8
2	B	568	VAL	3.8
2	B	543	GLU	3.7
3	L	161	LYS	3.7
2	B	467	PRO	3.7
2	B	517	GLY	3.6
2	B	542	ALA	3.6
2	B	569	GLU	3.6
2	B	280	VAL	3.5
2	B	571	LEU	3.5
2	B	279	PRO	3.4
3	L	157	SER	3.4
2	B	468	THR	3.3
2	B	554	GLN	3.3
4	H	139	LYS	3.2
2	B	506	LYS	3.1
4	H	201	THR	3.0
2	B	549	ARG	3.0
2	B	514	ALA	3.0
1	A	61	THR	3.0
2	B	289	LEU	2.9
4	H	200	GLY	2.9
2	B	504	GLY	2.8
2	B	551	HIS	2.8
2	B	505	GLY	2.6
2	B	516	GLY	2.6
2	B	548	THR	2.6

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Mol	Chain	Res	Type	RSRZ
3	L	1	GLU	2.6
1	A	151	ALA	2.6
2	B	492	SER	2.6
4	H	103	TRP	2.5
2	B	671	ALA	2.5
2	B	494	LYS	2.5
2	B	518	GLU	2.5
4	H	1	GLU	2.5
1	A	149	VAL	2.4
2	B	223	CYS	2.4
2	B	278	GLN	2.4
2	B	466	GLY	2.4
3	L	2	SER	2.4
4	H	137	SER	2.3
2	B	593	GLU	2.3
3	L	192	SER	2.3
4	H	167	GLY	2.3
2	B	314	ALA	2.2
2	B	591	HIS	2.2
2	B	179	LEU	2.2
3	L	62	ASP	2.2
2	B	303	ARG	2.2
2	B	672	VAL	2.2
2	B	539	ALA	2.2
2	B	658	SER	2.1
2	B	620	GLY	2.1
3	L	205	SER	2.1
3	L	154	LYS	2.1
2	B	499	ARG	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 carbohydrates in this entry.

6.4 Ligands

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95^{th} percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
5	NA	B	1	1/1	0.85	0.65	19.43	62,62,62,62	0

6.5 Other polymers

There are no such residues in this entry.