Full wwPDB X-ray Structure Validation Report (i)

Jan 31, 2016 – 07:35 PM GMT

PDB ID	:	1G9U
Title	:	CRYSTAL STRUCTURE OF YOPM-LEUCINE RICH EFFECTOR PROTEIN FROM YERSINIA PESTIS
Authors	:	Evdokimov, A.G.; Anderson, D.E.; Routzahn, K.M.; Waugh, D.S.
Deposited on	:	2000-11-28
Resolution	:	2.35 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:

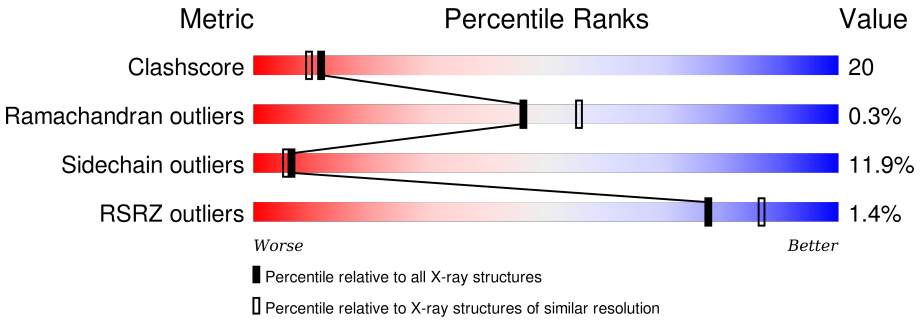
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.35 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))
Clashscore	102246	1456 (2.38-2.34)
Ramachandran outliers	100387	1435 (2.38-2.34)
Sidechain outliers	100360	1436 (2.38-2.34)
RSRZ outliers	91569	1358 (2.38-2.34)

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	454	% 43% 28% 6% 22%

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
2	ACT	A	1001	-	-	X	X
2	ACT	A	1002	-	-	-	X
2	ACT	A	1004	-	-	X	-
2	ACT	A	1005	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	ACT	A	1006	-	-	X	-
2	ACT	A	1007	-	-	-	X
2	ACT	A	1008	-	-	X	-
4	HG	A	2008	-	-	-	X

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 3075 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 OUTER PROTEIN YOPM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	353	Total	C	N	O	S	0	0	0
			2793	1773	457	558	5			

There are 45 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	385	ASN	-	SEE REMARK 999	UNP P17778
A	386	SER	-	SEE REMARK 999	UNP P17778
A	1387	HIS	-	SEE REMARK 999	UNP P17778
A	1388	LEU	-	SEE REMARK 999	UNP P17778
A	1389	ALA	-	SEE REMARK 999	UNP P17778
A	1390	GLU	-	SEE REMARK 999	UNP P17778
A	1391	VAL	-	SEE REMARK 999	UNP P17778
A	1392	PRO	-	SEE REMARK 999	UNP P17778
A	1393	GLU	-	SEE REMARK 999	UNP P17778
A	1394	LEU	-	SEE REMARK 999	UNP P17778
A	1395	PRO	-	SEE REMARK 999	UNP P17778
A	1396	GLN	-	SEE REMARK 999	UNP P17778
A	1397	ASN	-	SEE REMARK 999	UNP P17778
A	1398	LEU	-	SEE REMARK 999	UNP P17778
A	1399	LYS	-	SEE REMARK 999	UNP P17778
A	1400	GLN	-	SEE REMARK 999	UNP P17778
A	1401	LEU	-	SEE REMARK 999	UNP P17778
A	1402	HIS	-	SEE REMARK 999	UNP P17778
A	1403	VAL	-	SEE REMARK 999	UNP P17778
A	1404	GLU	-	SEE REMARK 999	UNP P17778
A	1405	THR	-	SEE REMARK 999	UNP P17778
A	1406	ASN	-	SEE REMARK 999	UNP P17778
A	1407	PRO	-	SEE REMARK 999	UNP P17778
A	1408	LEU	-	SEE REMARK 999	UNP P17778
A	1409	ARG	-	SEE REMARK 999	UNP P17778
A	1410	GLU	-	SEE REMARK 999	UNP P17778
A	1411	PHE	-	SEE REMARK 999	UNP P17778

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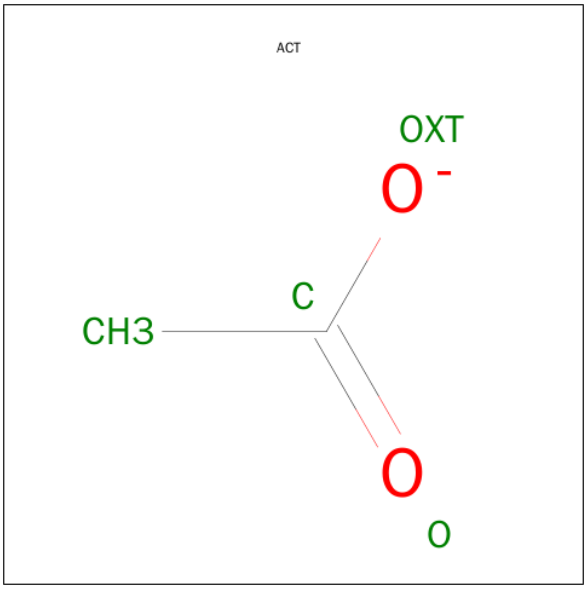
Chain	Residue	Modelled	Actual	Comment	Reference
A	1412	PRO	-	SEE REMARK 999	UNP P17778
A	1413	ASP	-	SEE REMARK 999	UNP P17778
A	1414	ILE	-	SEE REMARK 999	UNP P17778
A	1415	PRO	-	SEE REMARK 999	UNP P17778
A	1416	GLU	-	SEE REMARK 999	UNP P17778
A	1417	SER	-	SEE REMARK 999	UNP P17778
A	1418	VAL	-	SEE REMARK 999	UNP P17778
A	1419	GLU	-	SEE REMARK 999	UNP P17778
A	1420	ASP	-	SEE REMARK 999	UNP P17778
A	1421	LEU	-	SEE REMARK 999	UNP P17778
A	1422	ARG	-	SEE REMARK 999	UNP P17778
A	1423	MET	-	SEE REMARK 999	UNP P17778
A	1449	HIS	-	EXPRESSION TAG	UNP P17778
A	1450	HIS	-	EXPRESSION TAG	UNP P17778
A	1451	HIS	-	EXPRESSION TAG	UNP P17778
A	1452	HIS	-	EXPRESSION TAG	UNP P17778
A	1453	HIS	-	EXPRESSION TAG	UNP P17778
A	1454	HIS	-	EXPRESSION TAG	UNP P17778

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

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

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	2	Total Ca 2 2	0	0

- Molecule 4 is MERCURY (II) ION (three-letter code: HG) (formula: Hg).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	8	Total Hg 8 8	0	0

- Molecule 5 is water.

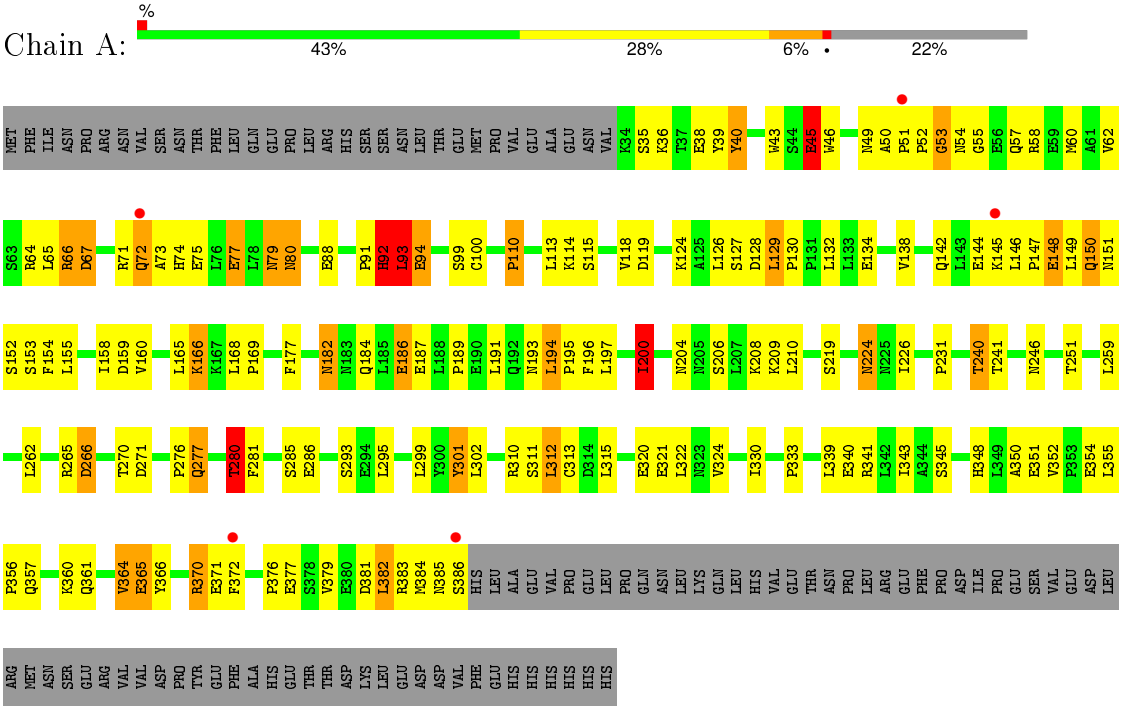
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	236	Total O 236 236	0	0

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: OUTER PROTEIN YOPM

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 42 2 2	Depositor
Cell constants a, b, c, α, β, γ	109.36Å 109.36Å 101.50Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	100.00 – 2.35 74.39 – 2.37	Depositor EDS
% Data completeness (in resolution range)	97.0 (100.00-2.35) 92.0 (74.39-2.37)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.13 (at 2.37Å)	Xtriage
Refinement program	SHELXL-97	Depositor
R, R_{free}	0.200 , 0.230 0.186 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å^2)	21.8	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 148.4	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Outliers	0 of 24660 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	3075	wwPDB-VP
Average B, all atoms (Å^2)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.21% 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: CA, HG, ACT

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	0.51	0/2853	1.64	39/3903 (1.0%)

There are no bond length outliers.

All (39) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	66	ARG	CD-NE-CZ	18.73	149.83	123.60
1	A	177	PHE	CB-CG-CD1	10.62	128.24	120.80
1	A	383	ARG	NE-CZ-NH1	8.00	124.30	120.30
1	A	40	TYR	CB-CG-CD1	-7.97	116.22	121.00
1	A	159	ASP	CB-CG-OD1	7.35	124.92	118.30
1	A	40	TYR	CB-CG-CD2	7.07	125.24	121.00
1	A	73	ALA	C-N-CA	7.05	139.32	121.70
1	A	301	TYR	CB-CG-CD2	7.03	125.22	121.00
1	A	313	CYS	CB-CA-C	6.94	124.29	110.40
1	A	64	ARG	NE-CZ-NH2	-6.61	116.99	120.30
1	A	64	ARG	NE-CZ-NH1	6.50	123.55	120.30
1	A	200	ILE	CB-CA-C	-6.42	98.75	111.60
1	A	383	ARG	NE-CZ-NH2	-6.37	117.11	120.30
1	A	53	GLY	C-N-CA	6.17	137.11	121.70
1	A	45	GLU	CB-CG-CD	6.05	130.55	114.20
1	A	312	LEU	C-N-CA	-6.03	106.61	121.70
1	A	128	ASP	CB-CG-OD1	5.98	123.69	118.30
1	A	310	ARG	NE-CZ-NH1	-5.93	117.33	120.30
1	A	92	HIS	C-N-CA	-5.91	106.93	121.70
1	A	311	SER	O-C-N	5.87	132.09	122.70
1	A	372	PHE	CA-CB-CG	5.85	127.94	113.90
1	A	66	ARG	NE-CZ-NH2	-5.81	117.39	120.30
1	A	266	ASP	CB-CG-OD2	5.79	123.51	118.30
1	A	93	LEU	CA-CB-CG	5.78	128.60	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	165	LEU	O-C-N	5.75	131.89	122.70
1	A	371	GLU	O-C-N	5.51	131.52	122.70
1	A	66	ARG	NE-CZ-NH1	5.50	123.05	120.30
1	A	80	ASN	N-CA-CB	-5.46	100.78	110.60
1	A	159	ASP	CB-CG-OD2	-5.43	113.41	118.30
1	A	72	GLN	CB-CG-CD	5.40	125.64	111.60
1	A	381	ASP	CB-CG-OD1	5.37	123.13	118.30
1	A	280	THR	N-CA-CB	-5.24	100.34	110.30
1	A	301	TYR	CB-CG-CD1	-5.18	117.89	121.00
1	A	128	ASP	CB-CG-OD2	-5.14	113.67	118.30
1	A	177	PHE	CB-CG-CD2	-5.13	117.21	120.80
1	A	382	LEU	CA-CB-CG	5.11	127.05	115.30
1	A	364	VAL	C-N-CA	5.09	134.42	121.70
1	A	45	GLU	OE1-CD-OE2	-5.04	117.26	123.30
1	A	153	SER	O-C-N	-5.01	114.68	122.70

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	2793	0	2781	110	0
2	A	36	0	27	12	0
3	A	2	0	0	0	0
4	A	8	0	0	1	0
5	A	236	0	0	13	0
All	All	3075	0	2808	114	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 20.

All (114) 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:92:HIS:CE1	4:A:2004:HG:HG	1.68	1.06
1:A:35:SER:OG	1:A:38:GLU:HG3	1.84	0.78
1:A:62:VAL:O	1:A:66:ARG:HG3	1.84	0.77
1:A:93:LEU:HD12	1:A:110:PRO:HG2	1.68	0.76
1:A:77:GLU:HG2	1:A:79:ASN:OD1	1.86	0.75
1:A:271:ASP:HB3	2:A:1003:ACT:H3	1.73	0.70
1:A:182:ASN:HD22	1:A:204:ASN:HB3	1.58	0.69
1:A:144:GLU:O	1:A:145:LYS:HD2	1.94	0.68
1:A:315:LEU:HD23	2:A:1008:ACT:H3	1.75	0.67
1:A:155:LEU:HD13	1:A:158:ILE:HD11	1.75	0.66
1:A:94:GLU:HG2	1:A:114:LYS:HD2	1.75	0.66
1:A:382:LEU:HD11	1:A:384:MET:HE2	1.77	0.65
1:A:130:PRO:HB2	5:A:3144:HOH:O	1.97	0.64
1:A:266:ASP:HA	1:A:286:GLU:O	1.98	0.63
1:A:376:PRO:HG2	1:A:379:VAL:HG23	1.78	0.63
1:A:80:ASN:HA	1:A:100:CYS:O	1.99	0.63
1:A:67:ASP:O	1:A:71:ARG:HG3	1.98	0.63
1:A:366:TYR:H	1:A:385:ASN:HB3	1.63	0.63
1:A:312:LEU:HB2	1:A:333:PRO:HD3	1.83	0.61
1:A:45:GLU:HG2	1:A:46:TRP:N	2.16	0.60
1:A:138:VAL:HG23	1:A:160:VAL:HG12	1.82	0.60
1:A:382:LEU:HD11	1:A:384:MET:CE	2.32	0.60
1:A:138:VAL:CG2	1:A:160:VAL:HG12	2.33	0.59
1:A:118:VAL:HG23	1:A:138:VAL:HG12	1.84	0.59
1:A:124:LYS:HE2	1:A:144:GLU:OE2	2.03	0.59
1:A:66:ARG:HD3	5:A:3130:HOH:O	2.02	0.58
1:A:147:PRO:HB2	2:A:1005:ACT:O	2.03	0.58
1:A:321:GLU:HG2	1:A:341:ARG:HB2	1.84	0.58
1:A:210:LEU:HB2	1:A:231:PRO:HD3	1.85	0.58
1:A:126:LEU:HD13	1:A:129:LEU:HD21	1.85	0.58
1:A:195:PRO:HG2	1:A:196:PHE:CD1	2.39	0.57
1:A:350:ALA:O	1:A:351:GLU:HG3	2.05	0.57
1:A:150:GLN:H	1:A:150:GLN:HE21	1.51	0.56
2:A:1005:ACT:H1	5:A:3218:HOH:O	2.06	0.56
1:A:193:ASN:HB2	5:A:3080:HOH:O	2.03	0.56
1:A:149:LEU:HD13	1:A:158:ILE:HD13	1.87	0.56
1:A:39:TYR:OH	1:A:91:PRO:HG3	2.06	0.56
1:A:52:PRO:HD2	5:A:3230:HOH:O	2.07	0.55
1:A:330:ILE:HG22	1:A:348:HIS:HB3	1.89	0.55
1:A:194:LEU:HD23	1:A:197:LEU:HD22	1.90	0.54
1:A:345:SER:OG	1:A:365:GLU:HG3	2.08	0.54
1:A:350:ALA:HB1	1:A:370:ARG:HG3	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:113:LEU:HB3	1:A:130:PRO:CG	2.38	0.53
1:A:352:VAL:CG1	1:A:355:LEU:HD21	2.38	0.53
1:A:49:ASN:HB3	2:A:1004:ACT:OXT	2.08	0.52
1:A:182:ASN:HD22	1:A:204:ASN:CB	2.21	0.52
2:A:1004:ACT:H1	5:A:3127:HOH:O	2.08	0.52
1:A:357:GLN:OE1	1:A:357:GLN:HA	2.08	0.52
1:A:186:GLU:OE1	1:A:208:LYS:HE2	2.09	0.52
1:A:50:ALA:HB2	1:A:58:ARG:HD2	1.92	0.51
1:A:197:LEU:HD21	1:A:200:ILE:HD13	1.93	0.51
1:A:341:ARG:HD2	5:A:3195:HOH:O	2.11	0.51
1:A:382:LEU:CD1	1:A:384:MET:HG3	2.41	0.50
1:A:351:GLU:HG3	1:A:370:ARG:HB2	1.93	0.50
1:A:149:LEU:HD12	1:A:169:PRO:HG2	1.93	0.50
1:A:382:LEU:HD11	1:A:384:MET:HG3	1.94	0.50
1:A:354:GLU:HA	5:A:3183:HOH:O	2.10	0.50
1:A:113:LEU:HB3	1:A:130:PRO:HG2	1.94	0.50
1:A:186:GLU:OE2	1:A:206:SER:O	2.30	0.50
1:A:151:ASN:HB2	2:A:1005:ACT:H2	1.93	0.49
1:A:67:ASP:OD2	1:A:71:ARG:HD2	2.12	0.49
1:A:281:PHE:CD1	1:A:301:TYR:HB3	2.47	0.49
1:A:118:VAL:CG2	1:A:138:VAL:HG12	2.43	0.49
1:A:295:LEU:HG	2:A:1008:ACT:H2	1.94	0.49
1:A:132:LEU:HB2	5:A:3144:HOH:O	2.14	0.48
1:A:343:ILE:HD11	1:A:361:GLN:OE1	2.13	0.48
1:A:364:VAL:O	1:A:364:VAL:HG22	2.14	0.48
1:A:99:SER:HB3	1:A:119:ASP:OD1	2.14	0.48
1:A:295:LEU:HD13	1:A:302:LEU:HD22	1.95	0.47
1:A:277:GLN:HG2	5:A:3207:HOH:O	2.14	0.47
1:A:240:THR:HG22	1:A:241:THR:OG1	2.14	0.47
1:A:182:ASN:ND2	1:A:204:ASN:HB3	2.29	0.47
1:A:191:LEU:HD13	1:A:200:ILE:CD1	2.45	0.47
1:A:352:VAL:HG11	1:A:355:LEU:HD21	1.97	0.47
1:A:146:LEU:HB3	1:A:147:PRO:HD2	1.96	0.47
1:A:166:LYS:HE2	1:A:184:GLN:OE1	2.15	0.47
1:A:145:LYS:HE3	5:A:3056:HOH:O	2.14	0.46
1:A:94:GLU:CG	1:A:114:LYS:HD2	2.43	0.46
1:A:280:THR:HA	1:A:299:LEU:HA	1.96	0.46
1:A:126:LEU:HA	1:A:126:LEU:HD23	1.79	0.46
1:A:384:MET:O	2:A:1001:ACT:H2	2.15	0.45
2:A:1006:ACT:H3	5:A:3047:HOH:O	2.16	0.45
1:A:265:ARG:HH11	1:A:265:ARG:HD2	1.66	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:365:GLU:H	1:A:365:GLU:HG3	1.37	0.45
1:A:129:LEU:HB3	1:A:130:PRO:HD2	1.98	0.45
1:A:134:GLU:HA	1:A:154:PHE:O	2.17	0.45
1:A:385:ASN:ND2	2:A:1001:ACT:O	2.49	0.45
1:A:148:GLU:OE1	1:A:150:GLN:NE2	2.50	0.45
1:A:57:GLN:HB3	1:A:60:MET:SD	2.56	0.45
1:A:312:LEU:HD23	1:A:312:LEU:HA	1.66	0.44
1:A:376:PRO:HG2	1:A:379:VAL:CG2	2.47	0.44
1:A:99:SER:O	1:A:100:CYS:HB2	2.18	0.44
1:A:355:LEU:HA	1:A:356:PRO:HD3	1.89	0.44
1:A:280:THR:HG23	1:A:281:PHE:CE2	2.52	0.44
1:A:322:LEU:HD11	1:A:324:VAL:HG13	1.99	0.44
2:A:1006:ACT:O	2:A:1009:ACT:H1	2.17	0.44
1:A:187:GLU:N	5:A:3209:HOH:O	2.50	0.44
1:A:36:LYS:HD3	1:A:40:TYR:OH	2.18	0.43
1:A:197:LEU:HD21	1:A:200:ILE:CD1	2.48	0.43
1:A:74:HIS:CD2	1:A:75:GLU:H	2.36	0.43
1:A:53:GLY:O	1:A:55:GLY:N	2.49	0.43
1:A:168:LEU:HB2	1:A:189:PRO:HD3	2.01	0.42
1:A:312:LEU:HD11	1:A:324:VAL:HG11	2.01	0.42
1:A:350:ALA:O	1:A:370:ARG:HB2	2.19	0.42
1:A:51:PRO:HA	1:A:52:PRO:HD3	1.58	0.42
1:A:265:ARG:HD2	1:A:285:SER:OG	2.20	0.42
1:A:224:ASN:HD22	1:A:246:ASN:HB3	1.85	0.41
1:A:364:VAL:HG13	1:A:364:VAL:O	2.21	0.41
1:A:259:LEU:HD21	1:A:262:LEU:HB2	2.02	0.41
1:A:43:TRP:CZ3	1:A:65:LEU:HB3	2.55	0.40
1:A:110:PRO:O	1:A:113:LEU:HB2	2.21	0.40
1:A:119:ASP:OD1	1:A:119:ASP:N	2.49	0.40
1:A:320:GLU:C	1:A:339:LEU:HD12	2.42	0.40
1:A:58:ARG:O	1:A:62:VAL:N	2.48	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	351/454 (77%)	314 (90%)	36 (10%)	1 (0%)	46 55

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	54	ASN

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	329/427 (77%)	290 (88%)	39 (12%)	6 5

All (39) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	45	GLU
1	A	67	ASP
1	A	72	GLN
1	A	77	GLU
1	A	79	ASN
1	A	88	GLU
1	A	92	HIS
1	A	93	LEU
1	A	94	GLU
1	A	110	PRO
1	A	115	SER
1	A	127	SER
1	A	129	LEU
1	A	142	GLN
1	A	148	GLU
1	A	150	GLN
1	A	152	SER
1	A	166	LYS

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Mol	Chain	Res	Type
1	A	182	ASN
1	A	186	GLU
1	A	194	LEU
1	A	200	ILE
1	A	209	LYS
1	A	219	SER
1	A	224	ASN
1	A	226	ILE
1	A	240	THR
1	A	251	THR
1	A	270	THR
1	A	276	PRO
1	A	277	GLN
1	A	280	THR
1	A	293	SER
1	A	340	GLU
1	A	360	LYS
1	A	365	GLU
1	A	370	ARG
1	A	377	GLU
1	A	386	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (9) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	41	ASN
1	A	74	HIS
1	A	80	ASN
1	A	122	ASN
1	A	150	GLN
1	A	182	ASN
1	A	192	GLN
1	A	224	ASN
1	A	358	ASN

5.3.3 RNA

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 carbohydrates in this entry.

5.6 Ligand geometry (i)

Of 19 ligands modelled in this entry, 10 are monoatomic - leaving 9 for Mogul analysis.

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	ACT	A	1001	-	1,3,3	0.05	0	0,3,3	0.00	-
2	ACT	A	1002	-	1,3,3	0.63	0	0,3,3	0.00	-
2	ACT	A	1003	-	1,3,3	0.92	0	0,3,3	0.00	-
2	ACT	A	1004	-	1,3,3	0.95	0	0,3,3	0.00	-
2	ACT	A	1005	-	1,3,3	0.49	0	0,3,3	0.00	-
2	ACT	A	1006	-	1,3,3	0.49	0	0,3,3	0.00	-
2	ACT	A	1007	-	1,3,3	0.06	0	0,3,3	0.00	-
2	ACT	A	1008	-	1,3,3	0.79	0	0,3,3	0.00	-
2	ACT	A	1009	-	1,3,3	0.74	0	0,3,3	0.00	-

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ACT	A	1001	-	-	0/0/0/0	0/0/0/0
2	ACT	A	1002	-	-	0/0/0/0	0/0/0/0
2	ACT	A	1003	-	-	0/0/0/0	0/0/0/0
2	ACT	A	1004	-	-	0/0/0/0	0/0/0/0
2	ACT	A	1005	-	-	0/0/0/0	0/0/0/0
2	ACT	A	1006	-	-	0/0/0/0	0/0/0/0

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	ACT	A	1007	-	-	0/0/0/0	0/0/0/0
2	ACT	A	1008	-	-	0/0/0/0	0/0/0/0
2	ACT	A	1009	-	-	0/0/0/0	0/0/0/0

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.

7 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1001	ACT	2	0
2	A	1003	ACT	1	0
2	A	1004	ACT	2	0
2	A	1005	ACT	3	0
2	A	1006	ACT	2	0
2	A	1008	ACT	2	0
2	A	1009	ACT	1	0

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 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	353/454 (77%)	-0.05	5 (1%)	78 87	12, 28, 67, 111	0

All (5) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	372	PHE	5.7
1	A	51	PRO	2.8
1	A	72	GLN	2.2
1	A	386	SER	2.2
1	A	145	LYS	2.2

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

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
4	HG	A	2008	1/1	0.93	0.36	42.03	92,92,92,92	1

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	ACT	A	1001	4/4	0.88	0.36	13.47	66,69,80,85	0
2	ACT	A	1002	4/4	0.95	0.20	8.22	16,51,55,61	0
2	ACT	A	1007	4/4	0.90	0.29	5.87	23,51,55,92	0
2	ACT	A	1006	4/4	0.94	0.16	1.56	37,43,53,68	0
2	ACT	A	1005	4/4	0.81	0.19	0.75	56,74,78,83	0
4	HG	A	2005	1/1	0.98	0.14	0.74	34,34,34,34	1
2	ACT	A	1004	4/4	0.81	0.21	0.56	51,70,72,75	0
3	CA	A	2001	1/1	0.99	0.11	0.34	20,20,20,20	0
2	ACT	A	1008	4/4	0.95	0.15	0.30	36,45,51,56	0
3	CA	A	2002	1/1	0.99	0.11	-1.02	22,22,22,22	0
2	ACT	A	1003	4/4	0.95	0.12	-1.89	39,44,49,49	0
4	HG	A	2003	1/1	0.99	0.07	-2.12	47,47,47,47	0
4	HG	A	2007	1/1	0.91	0.21	-	46,46,46,46	1
2	ACT	A	1009	4/4	0.86	0.30	-	80,84,87,88	0
4	HG	A	2004	1/1	0.97	0.10	-	29,29,29,29	1
4	HG	A	2009	1/1	0.54	0.26	-	96,96,96,96	1
4	HG	A	2006	1/1	0.57	0.31	-	112,112,112,112	1
4	HG	A	2010	1/1	0.75	0.42	-	157,157,157,157	1

6.5 Other polymers

There are no such residues in this entry.