Full wwPDB X-ray Structure Validation Report (i)

Feb 1, 2016 – 10:28 AM GMT

PDB ID	:	3M4Y
Title	:	Structural characterization of the subunit A mutant P235A of the A-ATP synthase
Authors	:	Manimekalai, M.S.; Balakrishna, A.M.; Kumar, A.; Priya, R.; Biukovic, G.; Jeyakanthan, J.; Gruber, G.
Deposited on	:	2010-03-12
Resolution	:	2.38 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:

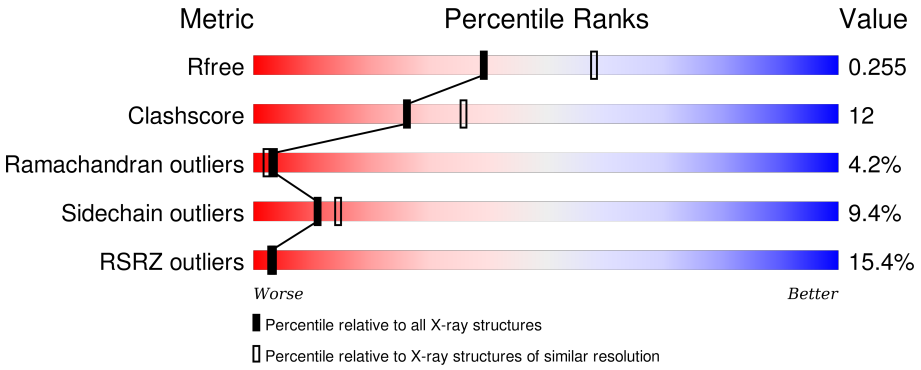
MolProbity	:	4.02b-467
Mogul	:	1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20026688
Percentile statistics	:	20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	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 2.38 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	4019 (2.40-2.36)
Clashscore	102246	4595 (2.40-2.36)
Ramachandran outliers	100387	4520 (2.40-2.36)
Sidechain outliers	100360	4522 (2.40-2.36)
RSRZ outliers	91569	4034 (2.40-2.36)

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	588	15% 71% 18% 6% • 5%

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	MPD	A	589	-	-	-	X

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MPD	A	590	-	-	-	X
2	MPD	A	591	-	-	X	-
3	TRS	A	592	-	-	-	X

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4899 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.

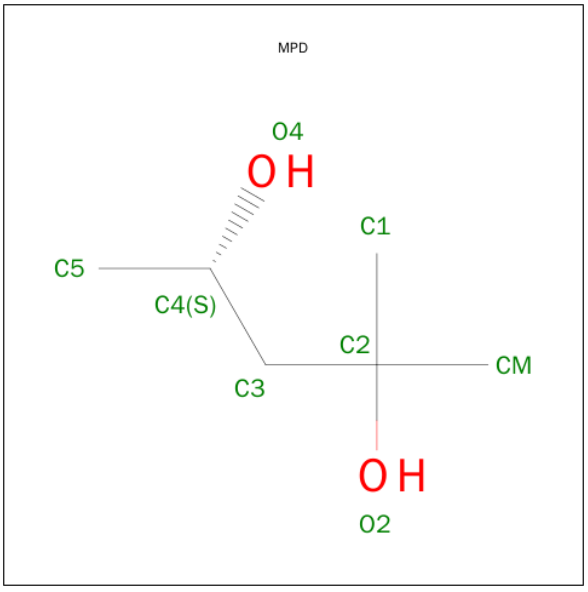
- Molecule 1 is a protein called V-type ATP synthase alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	558	Total	C	N	O	S	0	3	0
			4411	2821	752	820	18			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	235	ALA	PRO	ENGINEERED	UNP O57728

- Molecule 2 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



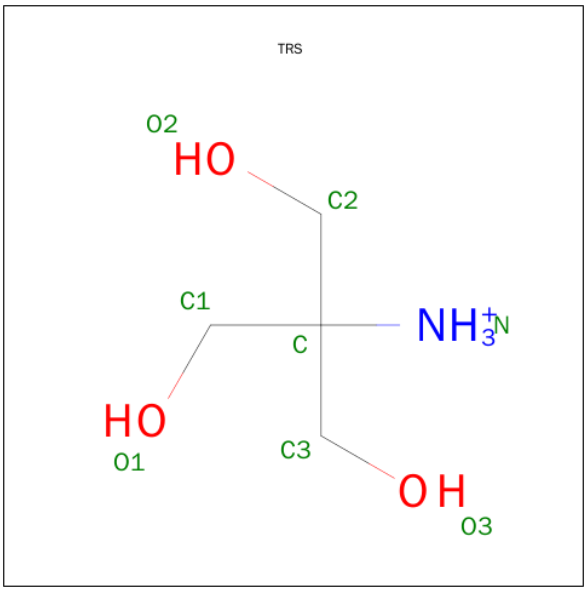
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			8	6	2		
2	A	1	Total	C	O	0	0
			8	6	2		

Continued on next page...

Continued from previous page...

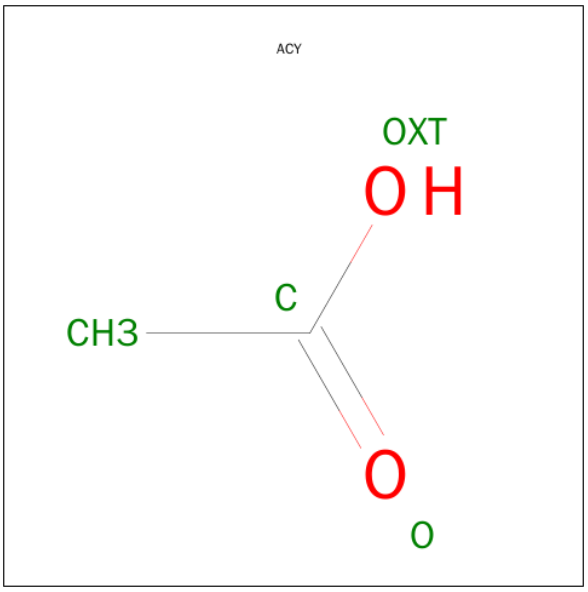
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 8 6 2	0	0
2	A	1	Total C O 8 6 2	0	0

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 8 4 1 3	0	0

- Molecule 4 is ACETIC ACID (three-letter code: ACY) (formula: $C_2H_4O_2$).



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

- Molecule 5 is water.

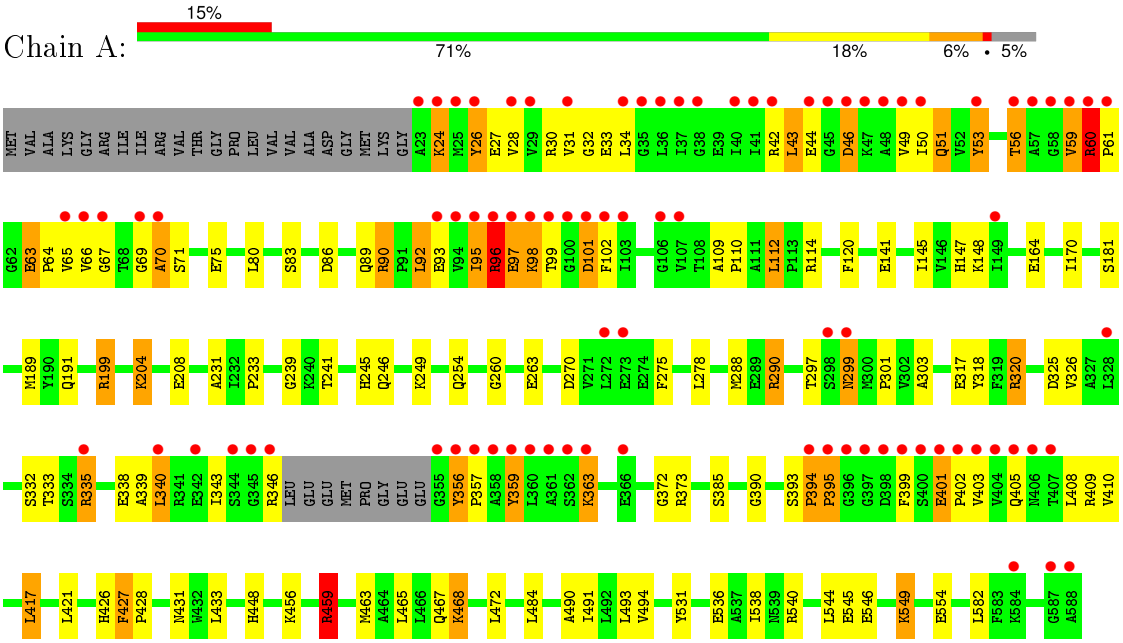
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	440	Total O 440 440	0	0

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: V-type ATP synthase alpha chain

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	127.56Å 127.56Å 105.40Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	24.93 – 2.38 24.93 – 2.38	Depositor EDS
% Data completeness (in resolution range)	99.8 (24.93-2.38) 99.8 (24.93-2.38)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	5.34 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.199 , 0.259 0.198 , 0.255	Depositor DCC
R_{free} test set	1769 reflections (5.27%)	DCC
Wilson B-factor (Å^2)	45.2	Xtriage
Anisotropy	0.046	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 67.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Outliers	0 of 35342 reflections	Xtriage
F_o,F_c correlation	0.94	EDS
Total number of atoms	4899	wwPDB-VP
Average B, all atoms (Å^2)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.08% 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: ACY, MPD, TRS

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	0.55	0/4512	0.67	2/6109 (0.0%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	427	PHE	C-N-CD	-11.29	95.75	120.60
1	A	459	ARG	CG-CD-NE	5.44	123.22	111.80

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	199	ARG	Sidechain
1	A	427	PHE	Peptide

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	4411	0	4480	105	0
2	A	32	0	56	11	0
3	A	8	0	12	1	0
4	A	8	0	6	0	0
5	A	440	0	0	5	0
All	All	4899	0	4554	111	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 12.

All (111) 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:394:PRO:HB2	1:A:395:PRO:HD3	1.14	1.13
1:A:394:PRO:HB2	1:A:395:PRO:CD	1.79	1.09
1:A:75:GLU:H	1:A:89:GLN:HE22	1.14	0.89
1:A:260:GLY:HA2	1:A:332:SER:HB3	1.54	0.89
1:A:120:PHE:HB2	1:A:189:MET:HE3	1.56	0.85
1:A:431:ASN:ND2	2:A:591:MPD:HM3	1.94	0.83
1:A:120:PHE:HB2	1:A:189:MET:CE	2.08	0.83
1:A:275:PHE:HB3	1:A:288[A]:MET:HE1	1.62	0.82
1:A:394:PRO:CB	1:A:395:PRO:HD3	2.05	0.79
1:A:199:ARG:NH1	1:A:318:TYR:HD1	1.83	0.76
1:A:356:TYR:HB3	1:A:359:TYR:HB3	1.69	0.75
1:A:405:GLN:HA	1:A:408:LEU:HB2	1.69	0.74
1:A:61:PRO:HD2	1:A:65:VAL:HG23	1.70	0.73
1:A:431:ASN:HD21	2:A:591:MPD:HM3	1.53	0.72
1:A:275:PHE:HB3	1:A:288[A]:MET:CE	2.19	0.71
1:A:401:GLU:H	1:A:402:PRO:HD3	1.55	0.70
1:A:199:ARG:HH11	1:A:318:TYR:HD1	1.38	0.69
1:A:86:ASP:OD1	1:A:90:ARG:HD3	1.94	0.68
2:A:595:MPD:O4	2:A:595:MPD:H11	1.93	0.68
1:A:44:GLU:HG2	1:A:53:TYR:CB	2.23	0.66
1:A:46:ASP:HB3	1:A:51:GLN:HB3	1.78	0.65
1:A:448:HIS:HE1	1:A:456:LYS:H	1.45	0.65
1:A:98:LYS:HG3	1:A:99:THR:H	1.62	0.64
1:A:141:GLU:OE1	1:A:147:HIS:HD2	1.80	0.63
1:A:549:LYS:C	1:A:549:LYS:HD2	2.19	0.63
2:A:591:MPD:HM2	2:A:591:MPD:O4	1.98	0.63
1:A:191:GLN:OE1	1:A:199:ARG:NH2	2.32	0.63

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:591:MPD:CM	2:A:591:MPD:O4	2.47	0.62
1:A:394:PRO:CB	1:A:395:PRO:CD	2.68	0.62
1:A:246:GLN:OE1	2:A:595:MPD:H13	1.99	0.61
1:A:401:GLU:N	1:A:402:PRO:CD	2.63	0.61
1:A:53:TYR:CD1	1:A:53:TYR:N	2.69	0.60
3:A:592:TRS:H21	5:A:713:HOH:O	2.00	0.60
1:A:494:VAL:HG11	1:A:531:TYR:HB2	1.82	0.60
2:A:591:MPD:H53	2:A:591:MPD:H11	1.84	0.59
1:A:53:TYR:HD1	1:A:53:TYR:N	2.00	0.59
1:A:59:VAL:HG23	1:A:67:GLY:H	1.68	0.59
1:A:96:ARG:HG3	1:A:97:GLU:H	1.69	0.58
1:A:239:GLY:HA2	5:A:633:HOH:O	2.03	0.58
1:A:204:LYS:HD2	1:A:372:GLY:HA3	1.84	0.58
1:A:24:LYS:HD2	1:A:26:TYR:H	1.68	0.58
1:A:468:LYS:HG3	1:A:491:ILE:HG21	1.86	0.58
1:A:44:GLU:HG2	1:A:53:TYR:HB2	1.85	0.58
1:A:69:GLY:O	1:A:70:ALA:HB2	2.04	0.56
1:A:463:MET:O	1:A:467:GLN:HG3	2.06	0.56
1:A:46:ASP:HB2	1:A:50:ILE:O	2.05	0.55
1:A:44:GLU:HG2	1:A:53:TYR:HB3	1.86	0.55
1:A:401:GLU:N	1:A:402:PRO:HD3	2.20	0.54
1:A:426:HIS:HE1	5:A:644:HOH:O	1.89	0.54
1:A:208:GLU:OE1	1:A:373:ARG:NH2	2.40	0.54
1:A:114:ARG:HD3	1:A:170:ILE:HD11	1.90	0.53
1:A:32:GLY:O	1:A:33:GLU:HB2	2.08	0.53
1:A:147:HIS:HE1	1:A:318:TYR:OH	1.92	0.53
1:A:546:GLU:HA	1:A:549:LYS:HE3	1.90	0.52
1:A:546:GLU:HG3	1:A:582:LEU:HD11	1.91	0.52
1:A:254:GLN:HE22	1:A:325:ASP:H	1.57	0.52
1:A:112:LEU:HB3	1:A:170:ILE:HD13	1.91	0.52
1:A:338:GLU:C	1:A:340:LEU:H	2.12	0.52
1:A:363:LYS:HB2	1:A:363:LYS:NZ	2.25	0.52
1:A:31:VAL:HB	1:A:301:PRO:HG3	1.90	0.52
1:A:95:ILE:HD13	1:A:96:ARG:N	2.26	0.51
1:A:408:LEU:C	1:A:410:VAL:H	2.15	0.50
1:A:260:GLY:CA	1:A:332:SER:HB3	2.34	0.50
1:A:59:VAL:O	1:A:60:ARG:HB2	2.11	0.50
1:A:490:ALA:HB2	1:A:538:ILE:CD1	2.42	0.50
1:A:109:ALA:HB1	1:A:110:PRO:CD	2.43	0.49
1:A:97:GLU:HA	1:A:101:ASP:O	2.13	0.47
1:A:545:GLU:H	1:A:545:GLU:CD	2.18	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:468:LYS:HG2	2:A:590:MPD:H13	1.97	0.46
1:A:317:GLU:O	1:A:320:ARG:HG3	2.15	0.46
1:A:63:GLU:O	1:A:65:VAL:N	2.49	0.46
1:A:26:TYR:CG	1:A:26:TYR:O	2.68	0.46
1:A:394:PRO:HG3	1:A:399:PHE:HB3	1.97	0.46
1:A:92:LEU:H	1:A:92:LEU:HD12	1.79	0.46
1:A:299:ASN:N	1:A:299:ASN:OD1	2.49	0.46
1:A:98:LYS:N	1:A:101:ASP:HA	2.30	0.45
1:A:335:ARG:N	1:A:335:ARG:HD2	2.31	0.45
1:A:233:PRO:HD2	1:A:417:LEU:HB2	1.98	0.45
1:A:148:LYS:NZ	5:A:668:HOH:O	2.49	0.45
1:A:98:LYS:HG3	1:A:99:THR:N	2.29	0.45
1:A:490:ALA:HB2	1:A:538:ILE:HD11	1.98	0.45
1:A:363:LYS:HB2	1:A:363:LYS:HZ2	1.83	0.44
1:A:249:LYS:NZ	2:A:595:MPD:H11	2.32	0.44
1:A:145:ILE:CG1	1:A:290:ARG:HB3	2.48	0.44
1:A:231:ALA:HA	1:A:390:GLY:O	2.17	0.44
1:A:333:THR:HG21	1:A:393:SER:H	1.83	0.44
1:A:254:GLN:NE2	1:A:325:ASP:H	2.15	0.43
1:A:408:LEU:O	1:A:410:VAL:N	2.51	0.43
2:A:591:MPD:H12	5:A:743:HOH:O	2.17	0.43
1:A:245[B]:HIS:CD2	2:A:595:MPD:H53	2.54	0.43
1:A:60:ARG:N	1:A:61:PRO:HD3	2.33	0.43
1:A:95:ILE:HD13	1:A:96:ARG:H	1.82	0.43
1:A:245[B]:HIS:CE1	1:A:278:LEU:HD13	2.53	0.43
1:A:549:LYS:HD2	1:A:549:LYS:O	2.18	0.43
1:A:373:ARG:HG3	1:A:385:SER:HB3	2.01	0.43
1:A:69:GLY:O	1:A:70:ALA:CB	2.67	0.42
1:A:33:GLU:HB3	1:A:303:ALA:HB3	2.01	0.42
1:A:50:ILE:HG22	1:A:66:VAL:HG13	2.02	0.42
1:A:317:GLU:HA	1:A:320:ARG:HG2	2.02	0.42
1:A:53:TYR:CE2	1:A:56:THR:O	2.72	0.42
1:A:468:LYS:HD2	1:A:472:LEU:HG	2.01	0.42
1:A:536:GLU:O	1:A:540:ARG:HG3	2.20	0.42
1:A:393:SER:HA	1:A:394:PRO:HD2	1.87	0.42
1:A:408:LEU:C	1:A:410:VAL:N	2.73	0.41
1:A:75:GLU:N	1:A:89:GLN:HE22	1.97	0.41
1:A:263:GLU:HA	1:A:335:ARG:HD3	2.01	0.41
1:A:544:LEU:HG	1:A:544:LEU:O	2.21	0.41
1:A:80:LEU:O	1:A:83:SER:HB2	2.20	0.41
1:A:459:ARG:HD3	1:A:459:ARG:HH11	1.71	0.41

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:356:TYR:HB3	1:A:359:TYR:CB	2.46	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	557/588 (95%)	505 (91%)	29 (5%)	23 (4%)	3 2

All (23) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	30	ARG
1	A	59	VAL
1	A	70	ALA
1	A	92	LEU
1	A	394	PRO
1	A	395	PRO
1	A	428	PRO
1	A	49	VAL
1	A	64	PRO
1	A	96	ARG
1	A	98	LYS
1	A	409	ARG
1	A	43	LEU
1	A	63	GLU
1	A	339	ALA
1	A	357	PRO
1	A	26	TYR
1	A	93	GLU
1	A	60	ARG
1	A	359	TYR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	356	TYR
1	A	403	VAL
1	A	401	GLU

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	471/491 (96%)	427 (91%)	44 (9%)	11 15

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	24	LYS
1	A	27	GLU
1	A	28	VAL
1	A	34	LEU
1	A	42	ARG
1	A	43	LEU
1	A	46	ASP
1	A	51	GLN
1	A	53	TYR
1	A	56	THR
1	A	60	ARG
1	A	71	SER
1	A	90	ARG
1	A	95	ILE
1	A	96	ARG
1	A	97	GLU
1	A	101	ASP
1	A	102	PHE
1	A	112	LEU
1	A	164	GLU
1	A	181	SER
1	A	204	LYS
1	A	241	THR

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	A	270	ASP
1	A	290	ARG
1	A	297	THR
1	A	299	ASN
1	A	320	ARG
1	A	326	VAL
1	A	335	ARG
1	A	340	LEU
1	A	343	ILE
1	A	346	ARG
1	A	363	LYS
1	A	417	LEU
1	A	421	LEU
1	A	433	LEU
1	A	459	ARG
1	A	465	LEU
1	A	468	LYS
1	A	484	LEU
1	A	493	LEU
1	A	549	LYS
1	A	554	GLU

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

Mol	Chain	Res	Type
1	A	89	GLN
1	A	147	HIS
1	A	254	GLN
1	A	426	HIS
1	A	448	HIS
1	A	504	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 carbohydrates in this entry.

5.6 Ligand geometry (i)

7 ligands are 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	MPD	A	589	-	6,7,7	0.28	0	7,10,10	0.38	0
2	MPD	A	590	-	6,7,7	0.24	0	7,10,10	0.50	0
2	MPD	A	591	-	6,7,7	0.29	0	7,10,10	0.16	0
3	TRS	A	592	-	7,7,7	0.97	1 (14%)	9,9,9	0.78	0
4	ACY	A	593	-	1,3,3	1.59	0	0,3,3	0.00	-
4	ACY	A	594	-	1,3,3	1.49	0	0,3,3	0.00	-
2	MPD	A	595	-	6,7,7	0.22	0	7,10,10	0.47	0

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	MPD	A	589	-	-	0/5/5/5	0/0/0/0
2	MPD	A	590	-	-	0/5/5/5	0/0/0/0
2	MPD	A	591	-	-	0/5/5/5	0/0/0/0
3	TRS	A	592	-	-	0/9/9/9	0/0/0/0
4	ACY	A	593	-	-	0/0/0/0	0/0/0/0
4	ACY	A	594	-	-	0/0/0/0	0/0/0/0
2	MPD	A	595	-	-	0/5/5/5	0/0/0/0

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	592	TRS	C-N	-2.47	1.47	1.50

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

4 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	590	MPD	1	0
2	A	591	MPD	6	0
3	A	592	TRS	1	0
2	A	595	MPD	4	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	558/588 (94%)	0.71	86 (15%) 3 3	24, 47, 123, 179	0

All (86) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	99	THR	11.4
1	A	26	TYR	9.6
1	A	397	GLY	8.7
1	A	399	PHE	8.4
1	A	358	ALA	8.1
1	A	23	ALA	7.7
1	A	398	ASP	7.6
1	A	403	VAL	7.3
1	A	402	PRO	7.2
1	A	357	PRO	6.7
1	A	407	THR	6.6
1	A	59	VAL	6.5
1	A	58	GLY	6.5
1	A	361	ALA	6.0
1	A	400	SER	5.8
1	A	396	GLY	5.7
1	A	100	GLY	5.6
1	A	101	ASP	5.6
1	A	41	ILE	5.5
1	A	401	GLU	5.4
1	A	395	PRO	5.2
1	A	356	TYR	5.2
1	A	103	ILE	5.0
1	A	344	SER	5.0
1	A	40	ILE	4.8
1	A	70	ALA	4.8
1	A	25	MET	4.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	98	LYS	4.6
1	A	102	PHE	4.4
1	A	355	GLY	4.4
1	A	404	VAL	4.4
1	A	107	VAL	4.3
1	A	405	GLN	4.3
1	A	65	VAL	4.1
1	A	95	ILE	4.1
1	A	46	ASP	4.1
1	A	94	VAL	4.1
1	A	60	ARG	4.0
1	A	69	GLY	3.9
1	A	50	ILE	3.8
1	A	406	ASN	3.8
1	A	359	TYR	3.7
1	A	35	GLY	3.6
1	A	298	SER	3.6
1	A	66	VAL	3.4
1	A	587	GLY	3.3
1	A	362	SER	3.2
1	A	49	VAL	3.2
1	A	394	PRO	3.2
1	A	24	LYS	3.2
1	A	106	GLY	3.1
1	A	67	GLY	3.0
1	A	31	VAL	2.9
1	A	96	ARG	2.9
1	A	273	GLU	2.9
1	A	36	LEU	2.8
1	A	366	GLU	2.8
1	A	345	GLY	2.8
1	A	93	GLU	2.8
1	A	584	LYS	2.8
1	A	346	ARG	2.7
1	A	342	GLU	2.7
1	A	34	LEU	2.7
1	A	47	LYS	2.7
1	A	48	ALA	2.6
1	A	42	ARG	2.6
1	A	37	ILE	2.6
1	A	149	ILE	2.5
1	A	44	GLU	2.5

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	53	TYR	2.5
1	A	97	GLU	2.4
1	A	340	LEU	2.4
1	A	29	VAL	2.4
1	A	45	GLY	2.4
1	A	360	LEU	2.3
1	A	28	VAL	2.3
1	A	56	THR	2.3
1	A	38	GLY	2.3
1	A	272	LEU	2.2
1	A	328	LEU	2.2
1	A	57	ALA	2.2
1	A	299	ASN	2.1
1	A	61	PRO	2.1
1	A	335	ARG	2.1
1	A	363	LYS	2.1
1	A	588	ALA	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 (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. 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
2	MPD	A	590	8/8	0.81	0.24	5.52	91,91,92,92	0
3	TRS	A	592	8/8	0.87	0.20	2.49	70,71,71,71	0
2	MPD	A	589	8/8	0.87	0.23	2.08	63,65,65,66	0
2	MPD	A	595	8/8	0.92	0.20	1.97	49,50,52,52	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å^2)	Q<0.9
2	MPD	A	591	8/8	0.91	0.20	0.58	62,66,67,67	0
4	ACY	A	594	4/4	0.93	0.13	0.43	72,72,72,72	0
4	ACY	A	593	4/4	0.76	0.36	-	74,74,74,74	0

6.5 Other polymers (i)

There are no such residues in this entry.