



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:46 PM GMT

PDB ID : 4AG5
Title : Structure of VirB4 of Thermoanaerobacter pseudethanolicus
Authors : Wallden, K.; Williams, R.; Yan, J.; Lian, P.W.; Wang, L.; Thalassinou, K.;
Orlova, E.V.; Waksman, G.
Deposited on : 2012-01-24
Resolution : 2.45 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 ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) 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

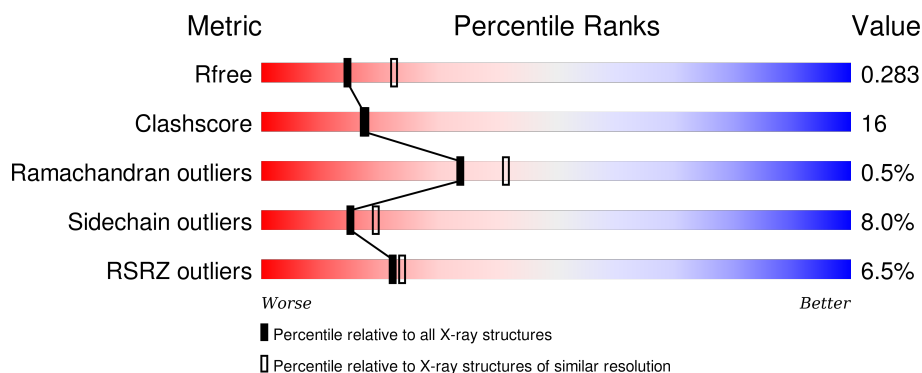
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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	4776 (2.50-2.42)
Clashscore	102246	1030 (2.48-2.44)
Ramachandran outliers	100387	1024 (2.48-2.44)
Sidechain outliers	100360	1024 (2.48-2.44)
RSRZ outliers	91569	4787 (2.50-2.42)

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 $\leq 5\%$. 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	392	<div> <div>68%</div> <div>24%</div> <div>• 5%</div> </div>
1	B	392	<div> <div>66%</div> <div>27%</div> <div>• 5%</div> </div>
1	C	392	<div> <div>5%</div> <div>46%</div> <div>20%</div> <div>• 32%</div> </div>
1	D	392	<div> <div>15%</div> <div>39%</div> <div>27%</div> <div>• 31%</div> </div>

2 Entry composition

There are 6 unique types of molecules in this entry. The entry contains 10083 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 TYPE IV SECRETORY PATHWAY VIRB4 COMPONENTS-LIKE PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	374	Total	C	N	O	S	0	0	0
			2941	1875	502	555	9			
1	B	374	Total	C	N	O	S	0	0	0
			2935	1868	502	556	9			
1	C	268	Total	C	N	O	S	0	0	0
			1994	1280	331	378	5			
1	D	270	Total	C	N	O	S	0	0	0
			1972	1265	327	377	3			

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		
2	C	1	Total	Mg	0	0
			1	1		

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			6	3	3		
3	A	1	Total	C	O	0	0
			6	3	3		
3	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
4	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

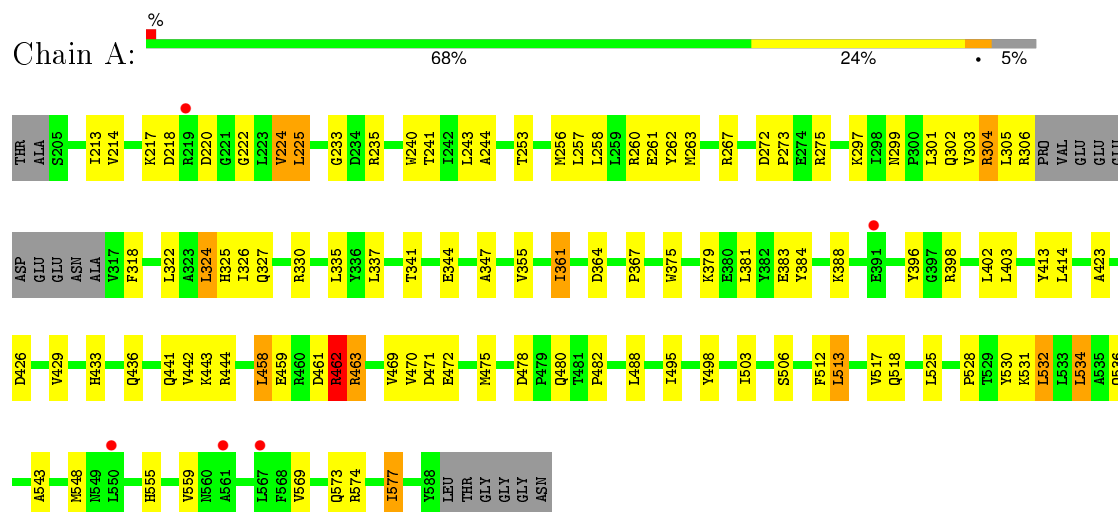
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	70	Total	O	0	0
			70	70		
6	B	65	Total	O	0	0
			65	65		
6	C	12	Total	O	0	0
			12	12		
6	D	14	Total	O	0	0
			14	14		

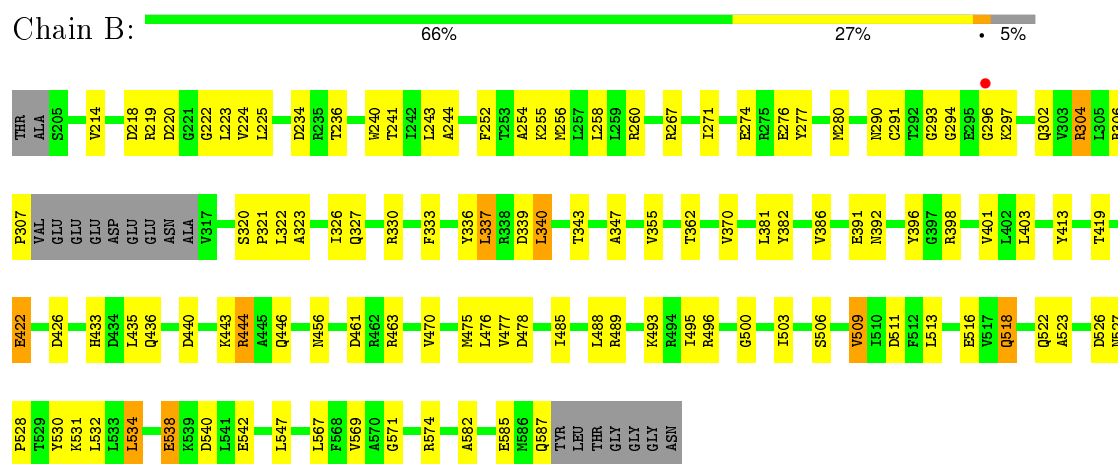
3 Residue-property plots

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.

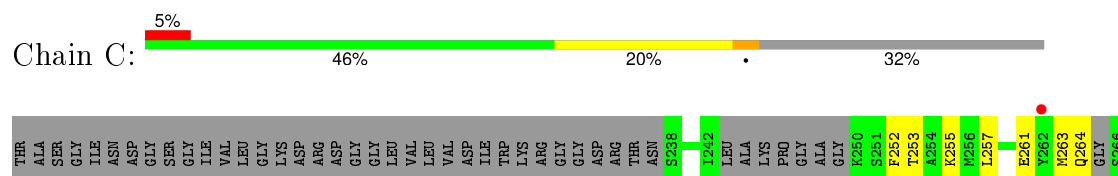
• Molecule 1: TYPE IV SECRETORY PATHWAY VIRB4 COMPONENTS-LIKE PROTEIN

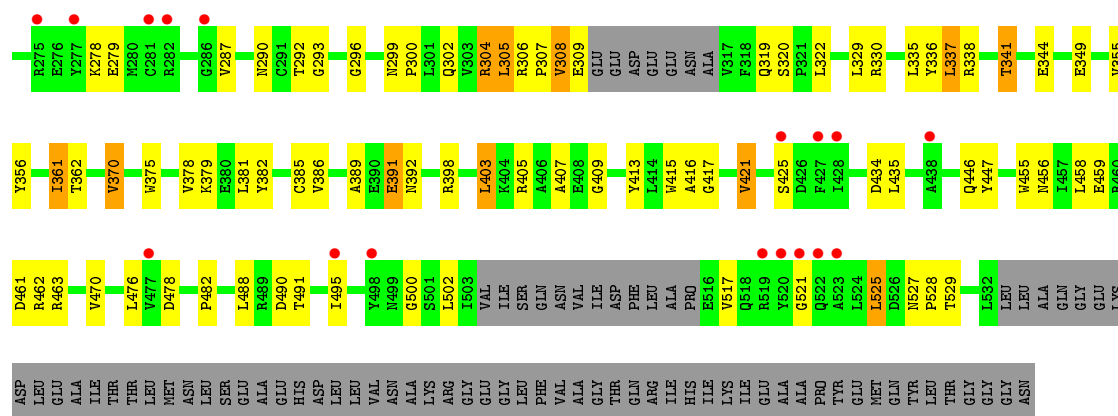


• Molecule 1: TYPE IV SECRETORY PATHWAY VIRB4 COMPONENTS-LIKE PROTEIN

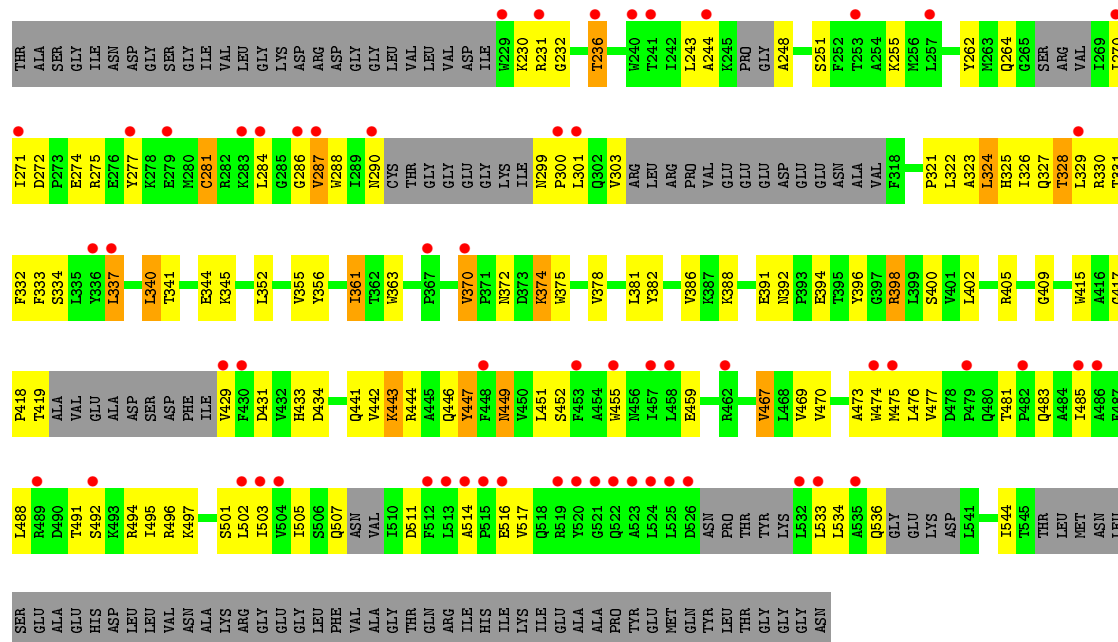


• Molecule 1: TYPE IV SECRETORY PATHWAY VIRB4 COMPONENTS-LIKE PROTEIN





● Molecule 1: TYPE IV SECRETORY PATHWAY VIRB4 COMPONENTS-LIKE PROTEIN



4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	108.08Å 110.80Å 156.43Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	54.04 – 2.45 54.04 – 2.45	Depositor EDS
% Data completeness (in resolution range)	95.8 (54.04-2.45) 95.4 (54.04-2.45)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.98 (at 2.45Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.233 , 0.288 0.227 , 0.283	Depositor DCC
R_{free} test set	3373 reflections (5.05%)	DCC
Wilson B-factor (Å ²)	46.2	Xtriage
Anisotropy	0.590	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 72.3	EDS
Estimated twinning fraction	0.023 for k,h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.30$	Xtriage
Outliers	1 of 69311 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10083	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.06% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

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

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	0.41	0/2998	0.56	0/4068
1	B	0.37	0/2992	0.56	0/4061
1	C	0.35	0/2036	0.55	0/2786
1	D	0.39	0/2005	0.61	3/2739 (0.1%)
All	All	0.38	0/10031	0.57	3/13654 (0.0%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	287	VAL	N-CA-CB	-7.53	94.94	111.50
1	D	287	VAL	N-CA-C	6.51	128.56	111.00
1	D	286	GLY	N-CA-C	-5.30	99.86	113.10

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	2941	0	2895	77	0
1	B	2935	0	2884	81	0
1	C	1994	0	1805	54	0
1	D	1972	0	1780	104	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	12	0	16	3	0
3	B	6	0	8	0	0
4	A	27	0	12	0	0
4	B	27	0	12	3	0
5	A	5	0	0	0	0
6	A	70	0	0	3	0
6	B	65	0	0	2	0
6	C	12	0	0	1	0
6	D	14	0	0	1	0
All	All	10083	0	9412	312	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 16.

All (312) 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:D:398:ARG:HH11	1:D:398:ARG:HG2	1.20	1.06
1:B:569:VAL:HG22	1:B:574:ARG:HG2	1.37	1.03
1:D:503:ILE:HG22	1:D:505:ILE:HD11	1.44	0.97
1:D:372:ASN:O	1:D:418:PRO:HB3	1.75	0.85
1:D:270:ILE:HD13	1:D:277:TYR:CD2	2.12	0.84
1:D:272:ASP:OD2	1:D:277:TYR:HD1	1.61	0.82
1:B:271:ILE:HD12	1:B:470:VAL:HG12	1.62	0.81
1:A:444:ARG:HG2	1:A:475:MET:O	1.81	0.81
1:D:449:ASN:HD22	1:D:449:ASN:C	1.86	0.79
1:C:341:THR:HG22	1:C:344:GLU:H	1.47	0.78
1:D:398:ARG:HH11	1:D:398:ARG:CG	1.93	0.78
1:D:270:ILE:CD1	1:D:277:TYR:CD2	2.68	0.77
1:C:338:ARG:HH12	1:C:478:ASP:CG	1.87	0.76
1:B:214:VAL:HG12	1:B:223:LEU:HD21	1.68	0.75
1:A:241:THR:HG23	1:A:528:PRO:HG2	1.68	0.75
1:A:218:ASP:HB2	1:A:222:GLY:H	1.52	0.75
1:D:391:GLU:HG3	1:D:392:ASN:HD22	1.53	0.74
1:B:509:VAL:HG13	1:B:540:ASP:CG	2.08	0.73
1:D:503:ILE:HG22	1:D:505:ILE:CD1	2.19	0.73
1:C:290:ASN:O	1:C:296:GLY:HA3	1.89	0.72
1:D:270:ILE:HD13	1:D:277:TYR:CE2	2.25	0.71

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:299:ASN:N	1:D:419:THR:HG1	1.88	0.71
1:C:382:TYR:O	1:C:386:VAL:HG23	1.92	0.70
1:C:287:VAL:HG12	1:C:425:SER:HB3	1.73	0.69
1:C:257:LEU:O	1:C:261:GLU:HG2	1.92	0.69
1:D:459:GLU:OE2	6:D:2011:HOH:O	2.09	0.69
1:C:305:LEU:HD23	1:C:459:GLU:HG2	1.75	0.69
1:A:555:HIS:O	1:A:559:VAL:HG23	1.93	0.69
1:A:573:GLN:HG3	3:A:1590:GOL:H11	1.77	0.67
1:D:323:ALA:HB2	1:D:363:TRP:HZ3	1.59	0.66
1:D:341:THR:O	1:D:345:LYS:HG3	1.95	0.66
1:C:361:ILE:HG13	1:C:361:ILE:O	1.96	0.65
1:D:503:ILE:CG2	1:D:505:ILE:HD11	2.22	0.65
1:D:329:LEU:O	1:D:333:PHE:HD2	1.78	0.65
4:B:1589:ADP:H4'	1:D:236:THR:HG21	1.78	0.65
1:B:339:ASP:CG	1:B:398:ARG:HH22	2.00	0.65
1:C:292:THR:HG21	1:C:434:ASP:HB2	1.78	0.65
1:D:398:ARG:NH1	1:D:398:ARG:HG2	2.00	0.64
1:D:447:TYR:OH	1:D:470:VAL:HG13	1.98	0.63
1:D:470:VAL:HG12	1:D:473:ALA:HB2	1.80	0.63
1:B:240:TRP:HB2	1:B:503:ILE:HG12	1.80	0.63
1:A:531:LYS:HB2	1:A:569:VAL:HB	1.79	0.63
1:D:467:VAL:HA	1:D:501:SER:O	1.99	0.63
1:D:455:TRP:CE3	1:D:491:THR:HG23	2.33	0.62
1:B:297:LYS:O	1:B:419:THR:HA	2.00	0.62
1:A:344:GLU:OE2	1:A:398:ARG:NH1	2.32	0.62
1:B:422:GLU:H	1:B:422:GLU:CD	2.02	0.62
1:D:321:PRO:O	1:D:324:LEU:HD12	2.00	0.62
1:B:321:PRO:HD2	6:B:2033:HOH:O	1.99	0.62
1:D:322:LEU:O	1:D:326:ILE:HG12	2.00	0.61
1:B:291:CYS:HA	1:B:296:GLY:O	1.98	0.61
1:A:218:ASP:HB3	1:A:220:ASP:H	1.64	0.61
1:A:426:ASP:HB3	1:A:463:ARG:HH22	1.66	0.61
1:B:355:VAL:HG21	1:B:381:LEU:HA	1.81	0.61
1:D:327:GLN:HA	1:D:330:ARG:HG2	1.80	0.61
1:C:527:ASN:N	1:C:528:PRO:HD2	2.15	0.61
1:D:337:LEU:HB3	1:D:340:LEU:HD13	1.83	0.60
1:C:252:PHE:O	1:C:255:LYS:HB2	2.01	0.60
1:D:441:GLN:HG2	1:D:442:VAL:N	2.16	0.60
1:C:336:TYR:CD2	1:C:337:LEU:HD13	2.37	0.60
1:B:516:GLU:N	1:B:516:GLU:OE1	2.28	0.60
1:D:270:ILE:HG12	1:D:469:VAL:HB	1.84	0.60

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:356:TYR:CD1	1:D:361:ILE:HD11	2.37	0.60
1:D:370:VAL:HG22	1:D:370:VAL:O	2.02	0.59
1:A:218:ASP:HB2	1:A:222:GLY:N	2.16	0.59
1:A:344:GLU:CD	1:A:398:ARG:HH12	2.06	0.59
1:C:461:ASP:O	1:C:463:ARG:N	2.32	0.59
1:A:305:LEU:HD23	1:A:367:PRO:HB2	1.84	0.59
1:B:382:TYR:O	1:B:386:VAL:HG23	2.03	0.58
1:B:241:THR:HG22	1:B:243:LEU:HG	1.84	0.58
1:B:538:GLU:CD	1:B:538:GLU:H	2.07	0.58
1:B:436:GLN:HA	1:B:443:LYS:NZ	2.19	0.57
1:D:324:LEU:HD12	1:D:325:HIS:H	1.69	0.57
1:C:385:CYS:O	1:C:389:ALA:HB2	2.04	0.57
1:A:384:TYR:CZ	1:A:388:LYS:NZ	2.68	0.57
1:D:323:ALA:HB2	1:D:363:TRP:CZ3	2.39	0.57
1:B:585:GLU:C	1:B:587:GLN:H	2.07	0.57
1:D:391:GLU:HG3	1:D:392:ASN:ND2	2.17	0.57
1:B:306:ARG:HB2	1:B:307:PRO:HD2	1.87	0.57
1:D:361:ILE:O	1:D:361:ILE:HG13	2.04	0.56
1:A:379:LYS:O	1:A:383:GLU:HG3	2.05	0.56
1:B:336:TYR:CD2	1:B:337:LEU:HD13	2.40	0.56
1:D:327:GLN:O	1:D:330:ARG:HG2	2.06	0.56
1:A:530:TYR:HA	1:A:569:VAL:O	2.05	0.56
1:B:337:LEU:HB3	1:B:340:LEU:HD13	1.86	0.56
1:A:241:THR:CG2	1:A:528:PRO:HG2	2.35	0.55
1:B:234:ASP:OD1	1:B:571:GLY:HA3	2.06	0.55
1:C:421:VAL:HA	6:C:2001:HOH:O	2.05	0.55
1:D:449:ASN:C	1:D:449:ASN:ND2	2.55	0.55
1:C:308:VAL:HG11	1:C:319:GLN:HE22	1.71	0.55
1:D:287:VAL:HG13	1:D:287:VAL:O	2.06	0.55
1:D:325:HIS:O	1:D:329:LEU:HD13	2.07	0.55
1:C:330:ARG:HD2	1:C:349:GLU:OE2	2.06	0.55
1:A:512:PHE:O	1:A:517:VAL:HG23	2.06	0.55
1:A:384:TYR:CE2	1:A:388:LYS:NZ	2.74	0.55
1:D:405:ARG:HA	1:D:409:GLY:HA3	1.88	0.55
1:A:573:GLN:CG	3:A:1590:GOL:H11	2.38	0.54
1:A:233:GLY:HA3	6:A:2007:HOH:O	2.05	0.54
1:A:461:ASP:O	1:A:463:ARG:HG2	2.07	0.54
1:D:321:PRO:HA	1:D:324:LEU:HD11	1.90	0.54
1:A:355:VAL:HG21	1:A:381:LEU:HA	1.89	0.54
1:D:441:GLN:HG2	1:D:442:VAL:H	1.72	0.54
1:A:224:VAL:HG11	1:A:577:ILE:HG13	1.89	0.54

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:526:ASP:HA	1:B:531:LYS:HE3	1.90	0.54
1:D:275:ARG:HE	1:D:431:ASP:CG	2.11	0.53
1:A:256:MET:O	1:A:260:ARG:HG3	2.08	0.53
1:A:569:VAL:HG22	1:A:574:ARG:HD3	1.90	0.53
1:D:270:ILE:CD1	1:D:277:TYR:CE2	2.90	0.53
1:D:446:GLN:OE1	1:D:446:GLN:HA	2.09	0.53
1:B:461:ASP:OD1	1:B:461:ASP:O	2.27	0.53
1:D:243:LEU:O	1:D:244:ALA:HB2	2.09	0.53
1:A:257:LEU:O	1:A:261:GLU:HG2	2.09	0.53
1:D:388:LYS:HB3	1:D:396:TYR:CD2	2.44	0.53
1:B:322:LEU:O	1:B:326:ILE:HG12	2.08	0.53
1:B:277:TYR:HD1	1:B:280:MET:CE	2.22	0.52
1:B:236:THR:O	1:B:496:ARG:HD3	2.09	0.52
1:C:379:LYS:HA	1:C:407:ALA:HB1	1.91	0.52
1:B:274:GLU:HB2	1:B:276:GLU:HG2	1.92	0.52
1:B:422:GLU:N	1:B:422:GLU:CD	2.62	0.52
1:A:235:ARG:HB3	6:A:2005:HOH:O	2.09	0.52
1:B:436:GLN:HA	1:B:443:LYS:HZ1	1.72	0.52
1:A:240:TRP:HB2	1:A:503:ILE:HG12	1.92	0.52
1:D:299:ASN:HB2	1:D:417:GLY:O	2.10	0.52
1:D:398:ARG:HH12	1:D:402:LEU:HD22	1.75	0.52
1:B:446:GLN:HA	1:B:446:GLN:OE1	2.10	0.51
1:C:370:VAL:HG22	1:C:375:TRP:CE2	2.45	0.51
1:C:527:ASN:N	1:C:528:PRO:CD	2.73	0.51
1:A:470:VAL:HG23	1:A:470:VAL:O	2.09	0.51
1:B:527:ASN:N	1:B:528:PRO:CD	2.73	0.51
1:A:213:ILE:HD11	1:A:257:LEU:HD12	1.92	0.51
1:D:355:VAL:HG21	1:D:381:LEU:HA	1.93	0.51
1:A:267:ARG:CZ	1:A:463:ARG:HG3	2.40	0.51
1:D:494:ARG:O	1:D:497:LYS:N	2.43	0.51
1:B:391:GLU:HG3	1:B:392:ASN:ND2	2.26	0.51
1:D:470:VAL:HG12	1:D:470:VAL:O	2.10	0.51
1:D:288:TRP:CD1	1:D:429:VAL:HB	2.45	0.51
1:A:347:ALA:HB1	1:A:396:TYR:CZ	2.47	0.50
1:A:513:LEU:HD23	1:A:543:ALA:HB1	1.94	0.50
1:D:443:LYS:HD2	1:D:444:ARG:N	2.26	0.50
1:B:254:ALA:O	1:B:258:LEU:HG	2.12	0.50
1:B:426:ASP:CG	1:B:463:ARG:HH22	2.14	0.50
1:A:305:LEU:CD2	1:A:367:PRO:HB2	2.41	0.50
1:A:573:GLN:CD	3:A:1590:GOL:H11	2.32	0.50
1:A:301:LEU:HD23	1:A:301:LEU:N	2.27	0.50

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:322:LEU:O	1:A:326:ILE:HG12	2.12	0.50
1:B:290:ASN:O	1:B:296:GLY:HA3	2.12	0.49
1:D:325:HIS:HA	1:D:328:THR:HG23	1.95	0.49
1:B:433:HIS:C	1:B:435:LEU:H	2.15	0.49
1:C:521:GLY:O	1:C:525:LEU:HG	2.11	0.49
1:D:511:ASP:O	1:D:517:VAL:HG21	2.12	0.49
1:A:398:ARG:HB3	1:A:398:ARG:HH11	1.78	0.49
1:D:275:ARG:HG2	1:D:431:ASP:OD2	2.12	0.49
1:B:304:ARG:HB2	1:B:456:ASN:CG	2.31	0.49
1:A:241:THR:HG21	1:A:525:LEU:HA	1.94	0.49
1:D:303:VAL:HA	1:D:325:HIS:CG	2.48	0.49
1:B:413:TYR:OH	1:B:435:LEU:HD21	2.13	0.49
1:B:347:ALA:HB1	1:B:396:TYR:CZ	2.47	0.49
1:A:302:GLN:O	1:A:304:ARG:HD3	2.13	0.49
1:A:214:VAL:O	1:A:260:ARG:NH1	2.45	0.49
1:B:256:MET:HB3	1:B:260:ARG:HH11	1.78	0.49
1:C:461:ASP:OD1	1:C:461:ASP:O	2.30	0.49
1:C:308:VAL:HG11	1:C:319:GLN:NE2	2.28	0.49
1:D:477:VAL:HA	1:D:485:ILE:HD11	1.95	0.49
1:A:262:TYR:HD2	1:A:263:MET:HE2	1.78	0.49
1:A:444:ARG:HD2	1:A:478:ASP:HB2	1.95	0.48
1:C:344:GLU:OE1	1:C:398:ARG:NH1	2.46	0.48
1:C:308:VAL:HG22	1:C:309:GLU:N	2.27	0.48
1:C:391:GLU:CD	1:C:392:ASN:HD22	2.16	0.48
1:A:518:GLN:NE2	1:B:542:GLU:HG3	2.27	0.48
1:A:433:HIS:O	1:A:436:GLN:HG2	2.14	0.48
1:B:585:GLU:C	1:B:587:GLN:N	2.67	0.48
1:C:455:TRP:CZ3	1:C:458:LEU:HD23	2.47	0.48
1:B:255:LYS:HG2	1:B:280:MET:HE2	1.95	0.48
1:D:274:GLU:HA	1:D:433:HIS:ND1	2.28	0.48
1:B:477:VAL:HA	1:B:485:ILE:HD11	1.95	0.48
1:A:299:ASN:C	1:A:299:ASN:OD1	2.52	0.48
1:B:398:ARG:HA	1:B:401:VAL:HG12	1.96	0.47
1:D:251:SER:O	1:D:255:LYS:HG2	2.13	0.47
1:A:213:ILE:CD1	1:A:257:LEU:HD12	2.44	0.47
1:B:444:ARG:HD3	1:B:475:MET:O	2.14	0.47
1:B:218:ASP:HB3	1:B:220:ASP:H	1.79	0.47
1:D:514:ALA:HB3	1:D:517:VAL:HG22	1.97	0.47
1:C:304:ARG:HB2	1:C:456:ASN:CG	2.35	0.47
1:B:530:TYR:HA	1:B:569:VAL:O	2.15	0.47
1:B:478:ASP:OD1	1:B:478:ASP:C	2.52	0.47

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:327:GLN:O	1:B:330:ARG:HB2	2.15	0.47
1:B:267:ARG:NH1	1:B:463:ARG:HD3	2.29	0.47
1:D:476:LEU:HD13	1:D:485:ILE:HD13	1.97	0.47
1:D:290:ASN:HA	1:D:431:ASP:HB3	1.97	0.47
1:D:374:LYS:HG3	1:D:374:LYS:O	2.14	0.47
1:B:333:PHE:O	1:B:337:LEU:HB2	2.15	0.47
1:D:492:SER:HB2	1:D:502:LEU:HD22	1.97	0.47
1:A:263:MET:CE	1:A:263:MET:HA	2.45	0.46
1:D:443:LYS:C	1:D:443:LYS:HD2	2.36	0.46
1:D:444:ARG:HG3	1:D:475:MET:O	2.15	0.46
1:C:446:GLN:OE1	1:C:446:GLN:HA	2.14	0.46
1:B:243:LEU:O	1:B:244:ALA:HB2	2.15	0.46
1:A:534:LEU:O	1:A:536:GLN:NE2	2.49	0.46
1:B:582:ALA:O	1:B:585:GLU:HB2	2.15	0.46
1:B:523:ALA:O	1:B:527:ASN:HB2	2.15	0.46
1:C:403:LEU:HD12	1:C:403:LEU:HA	1.79	0.46
1:A:306:ARG:C	1:A:459:GLU:OE1	2.54	0.46
1:A:459:GLU:HG2	1:A:498:TYR:OH	2.15	0.46
1:D:244:ALA:O	1:D:507:GLN:HA	2.16	0.46
1:D:470:VAL:CG1	1:D:473:ALA:HB2	2.44	0.46
1:D:495:ILE:HD13	1:D:501:SER:C	2.37	0.46
1:C:491:THR:HG22	1:C:502:LEU:HD22	1.98	0.46
1:B:538:GLU:CD	1:B:538:GLU:N	2.69	0.45
1:D:248:ALA:O	1:D:534:LEU:HD23	2.16	0.45
1:B:219:ARG:HG3	1:B:220:ASP:OD1	2.16	0.45
1:D:262:TYR:C	1:D:264:GLN:N	2.69	0.45
1:A:361:ILE:HD13	1:A:375:TRP:HZ3	1.81	0.45
1:D:324:LEU:HD12	1:D:325:HIS:N	2.31	0.45
1:D:331:THR:O	1:D:332:PHE:C	2.52	0.45
1:A:272:ASP:OD1	1:A:275:ARG:N	2.49	0.45
1:B:436:GLN:HA	1:B:436:GLN:OE1	2.17	0.45
1:A:482:PRO:HB3	1:A:517:VAL:HG12	1.99	0.45
1:C:338:ARG:NH1	1:C:478:ASP:OD2	2.48	0.45
1:C:488:LEU:O	1:C:491:THR:HB	2.17	0.45
1:A:258:LEU:HD11	1:A:469:VAL:HG21	1.98	0.45
1:D:481:THR:C	1:D:483:GLN:H	2.19	0.45
1:A:304:ARG:HG2	1:A:325:HIS:CE1	2.52	0.45
1:A:337:LEU:HD22	1:A:402:LEU:HB3	1.99	0.45
1:C:306:ARG:HB2	1:C:307:PRO:HD2	1.99	0.45
1:C:263:MET:O	1:C:264:GLN:C	2.55	0.44
1:C:322:LEU:HD11	1:C:356:TYR:CD2	2.52	0.44

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:213:ILE:O	1:A:225:LEU:HD23	2.17	0.44
1:D:536:GLN:NE2	1:D:544:ILE:HD13	2.33	0.44
1:B:256:MET:HB3	1:B:260:ARG:NH1	2.33	0.44
1:D:329:LEU:HD23	1:D:352:LEU:HD13	1.99	0.44
1:C:308:VAL:HG21	1:C:319:GLN:OE1	2.17	0.44
1:A:327:GLN:O	1:A:330:ARG:HB2	2.17	0.44
1:B:440:ASP:O	1:B:443:LYS:HB2	2.16	0.44
1:D:382:TYR:O	1:D:386:VAL:HG23	2.17	0.44
1:B:320:SER:HB3	1:B:323:ALA:HB3	1.99	0.44
1:B:343:THR:HG23	1:C:320:SER:HB3	1.99	0.44
1:B:252:PHE:CE1	4:B:1589:ADP:C6	3.06	0.44
1:D:271:ILE:HD11	1:D:451:LEU:HA	1.99	0.44
1:D:533:LEU:HD22	1:D:544:ILE:HD12	1.99	0.44
1:A:459:GLU:HG2	1:A:498:TYR:CE2	2.52	0.43
1:B:255:LYS:HG2	1:B:280:MET:CE	2.47	0.43
1:C:302:GLN:HG3	1:C:375:TRP:CD1	2.53	0.43
1:B:477:VAL:HG11	1:B:511:ASP:O	2.17	0.43
1:B:522:GLN:HG3	1:B:547:LEU:HD11	2.00	0.43
1:C:300:PRO:O	1:C:329:LEU:HD21	2.19	0.43
1:D:398:ARG:NH1	1:D:398:ARG:CG	2.62	0.43
1:B:297:LYS:HE2	1:B:413:TYR:O	2.18	0.43
1:A:253:THR:HG21	1:A:532:LEU:CD1	2.49	0.43
1:D:495:ILE:CG2	1:D:496:ARG:N	2.81	0.43
1:B:477:VAL:HG22	1:B:485:ILE:CD1	2.49	0.43
1:D:262:TYR:O	1:D:264:GLN:N	2.52	0.43
1:A:532:LEU:HD23	1:A:532:LEU:HA	1.93	0.43
1:A:442:VAL:O	1:A:443:LYS:C	2.58	0.42
1:A:458:LEU:HD22	1:A:495:ILE:HG13	2.00	0.42
1:D:262:TYR:C	1:D:264:GLN:H	2.21	0.42
1:A:324:LEU:HA	1:A:324:LEU:HD12	1.86	0.42
4:B:1589:ADP:H4'	1:D:236:THR:CG2	2.48	0.42
1:B:218:ASP:HB2	1:B:222:GLY:N	2.34	0.42
1:D:382:TYR:HE1	1:D:400:SER:O	2.01	0.42
1:C:300:PRO:CD	1:C:415:TRP:HB3	2.50	0.42
1:C:293:GLY:HA2	1:C:413:TYR:O	2.20	0.42
1:B:294:GLY:HA2	1:B:297:LYS:HE3	2.02	0.42
1:D:303:VAL:HB	1:D:375:TRP:CH2	2.55	0.42
1:C:306:ARG:HB2	1:C:307:PRO:CD	2.50	0.42
1:D:327:GLN:O	1:D:328:THR:C	2.57	0.42
1:C:278:LYS:O	1:C:279:GLU:C	2.58	0.42
1:A:413:TYR:CE2	1:A:414:LEU:CD2	3.03	0.42

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:567:LEU:CD2	1:B:574:ARG:HD2	2.50	0.41
1:A:531:LYS:HD3	1:A:548:MET:CE	2.50	0.41
1:C:287:VAL:CG1	1:C:425:SER:HB3	2.45	0.41
1:A:303:VAL:HG13	1:A:325:HIS:HB2	2.02	0.41
1:B:337:LEU:HD12	1:B:337:LEU:HA	1.74	0.41
1:C:335:LEU:HA	1:C:335:LEU:HD23	1.85	0.41
1:C:495:ILE:HG12	1:C:500:GLY:HA3	2.03	0.41
1:A:335:LEU:O	1:A:441:GLN:HB2	2.20	0.41
1:B:477:VAL:HG22	1:B:485:ILE:HD11	2.01	0.41
1:A:471:ASP:CG	6:A:2017:HOH:O	2.59	0.41
1:B:495:ILE:HG12	1:B:500:GLY:HA3	2.02	0.41
1:D:275:ARG:NH2	1:D:434:ASP:OD1	2.52	0.41
1:D:281:CYS:SG	1:D:288:TRP:HB2	2.60	0.41
1:C:307:PRO:HB2	1:C:490:ASP:HB3	2.01	0.41
1:D:270:ILE:HG12	1:D:469:VAL:CG1	2.50	0.41
1:B:214:VAL:HG12	1:B:223:LEU:CD2	2.46	0.41
1:B:509:VAL:HG13	1:B:540:ASP:CB	2.49	0.41
1:D:474:TRP:HA	1:D:477:VAL:HG23	2.03	0.41
1:A:273:PRO:HG2	1:A:472:GLU:HB2	2.02	0.41
1:D:270:ILE:HD13	1:D:277:TYR:CG	2.54	0.41
1:D:303:VAL:HG13	1:D:325:HIS:HB2	2.03	0.41
1:A:303:VAL:HG13	1:A:325:HIS:CB	2.51	0.41
1:A:243:LEU:O	1:A:244:ALA:HB2	2.20	0.41
1:D:341:THR:OG1	1:D:344:GLU:HG3	2.21	0.41
1:D:287:VAL:O	1:D:287:VAL:CG1	2.69	0.41
1:B:496:ARG:NH2	6:B:2056:HOH:O	2.54	0.41
1:C:355:VAL:HG21	1:C:381:LEU:HA	2.03	0.41
1:D:300:PRO:CD	1:D:415:TRP:HB3	2.51	0.41
1:A:364:ASP:CG	1:D:394:GLU:HB2	2.41	0.41
1:B:518:GLN:HE21	1:B:518:GLN:HB2	1.70	0.41
1:C:482:PRO:HB3	1:C:517:VAL:HG12	2.03	0.41
1:A:318:PHE:CZ	1:D:394:GLU:HB3	2.55	0.41
1:B:302:GLN:O	1:B:304:ARG:HD3	2.21	0.40
1:A:462:ARG:H	1:A:462:ARG:HG2	1.51	0.40
1:D:262:TYR:CD2	1:D:262:TYR:C	2.94	0.40
1:C:413:TYR:OH	1:C:435:LEU:HD21	2.21	0.40
1:C:476:LEU:HD12	1:C:476:LEU:HA	1.83	0.40
1:C:299:ASN:HB2	1:C:417:GLY:O	2.20	0.40
1:C:405:ARG:HA	1:C:409:GLY:HA3	2.02	0.40
1:B:214:VAL:CG1	1:B:223:LEU:HD21	2.42	0.40
1:D:481:THR:O	1:D:481:THR:HG22	2.21	0.40

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:231:ARG:HG3	1:D:232:GLY:N	2.37	0.40
1:D:270:ILE:O	1:D:270:ILE:HG22	2.21	0.40
1:A:297:LYS:NZ	1:A:297:LYS:HB2	2.36	0.40
1:B:244:ALA:HB2	1:B:534:LEU:HB2	2.03	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	370/392 (94%)	351 (95%)	17 (5%)	2 (0%)	34	41
1	B	370/392 (94%)	346 (94%)	23 (6%)	1 (0%)	46	57
1	C	258/392 (66%)	228 (88%)	28 (11%)	2 (1%)	24	28
1	D	252/392 (64%)	239 (95%)	12 (5%)	1 (0%)	39	49
All	All	1250/1568 (80%)	1164 (93%)	80 (6%)	6 (0%)	34	41

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	462	ARG
1	B	293	GLY
1	C	462	ARG
1	D	230	LYS
1	A	423	ALA
1	C	416	ALA

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	304/330 (92%)	285 (94%)	19 (6%)	22	30
1	B	304/330 (92%)	283 (93%)	21 (7%)	19	25
1	C	186/330 (56%)	169 (91%)	17 (9%)	12	13
1	D	179/330 (54%)	158 (88%)	21 (12%)	7	6
All	All	973/1320 (74%)	895 (92%)	78 (8%)	15	19

All (78) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	217	LYS
1	A	224	VAL
1	A	225	LEU
1	A	304	ARG
1	A	324	LEU
1	A	341	THR
1	A	361	ILE
1	A	403	LEU
1	A	429	VAL
1	A	458	LEU
1	A	462	ARG
1	A	463	ARG
1	A	480	GLN
1	A	488	LEU
1	A	506	SER
1	A	513	LEU
1	A	532	LEU
1	A	534	LEU
1	A	577	ILE
1	B	224	VAL
1	B	225	LEU
1	B	304	ARG
1	B	337	LEU
1	B	340	LEU
1	B	362	THR
1	B	370	VAL
1	B	403	LEU
1	B	422	GLU

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	B	444	ARG
1	B	476	LEU
1	B	488	LEU
1	B	489	ARG
1	B	493	LYS
1	B	506	SER
1	B	509	VAL
1	B	513	LEU
1	B	518	GLN
1	B	532	LEU
1	B	534	LEU
1	B	538	GLU
1	C	253	THR
1	C	304	ARG
1	C	305	LEU
1	C	308	VAL
1	C	337	LEU
1	C	341	THR
1	C	361	ILE
1	C	362	THR
1	C	370	VAL
1	C	378	VAL
1	C	391	GLU
1	C	403	LEU
1	C	421	VAL
1	C	447	TYR
1	C	470	VAL
1	C	525	LEU
1	C	529	THR
1	D	236	THR
1	D	281	CYS
1	D	284	LEU
1	D	301	LEU
1	D	324	LEU
1	D	328	THR
1	D	334	SER
1	D	337	LEU
1	D	340	LEU
1	D	361	ILE
1	D	370	VAL
1	D	374	LYS
1	D	378	VAL

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type
1	D	398	ARG
1	D	443	LYS
1	D	447	TYR
1	D	449	ASN
1	D	452	SER
1	D	467	VAL
1	D	488	LEU
1	D	516	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	208	ASN
1	A	290	ASN
1	A	319	GLN
1	A	518	GLN
1	A	555	HIS
1	B	518	GLN
1	B	527	ASN
1	D	239	ASN
1	D	480	GLN
1	D	536	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](#)

Of 9 ligands modelled in this entry, 3 are monoatomic - leaving 6 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$
3	GOL	A	1590	-	5,5,5	0.33	0	5,5,5	0.35	0
4	ADP	A	1591	2	22,29,29	1.05	2 (9%)	27,45,45	1.75	3 (11%)
5	SO4	A	1592	-	4,4,4	0.18	0	6,6,6	0.08	0
3	GOL	A	1593	-	5,5,5	0.31	0	5,5,5	0.23	0
4	ADP	B	1589	2	22,29,29	1.11	2 (9%)	27,45,45	1.68	3 (11%)
3	GOL	B	1590	-	5,5,5	0.37	0	5,5,5	0.26	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
3	GOL	A	1590	-	-	0/4/4/4	0/0/0/0
4	ADP	A	1591	2	-	0/12/32/32	0/3/3/3
5	SO4	A	1592	-	-	0/0/0/0	0/0/0/0
3	GOL	A	1593	-	-	0/4/4/4	0/0/0/0
4	ADP	B	1589	2	-	0/12/32/32	0/3/3/3
3	GOL	B	1590	-	-	0/4/4/4	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1589	ADP	O4'-C1'	2.13	1.43	1.41
4	A	1591	ADP	O4'-C1'	2.37	1.44	1.41
4	A	1591	ADP	C5-C4	3.01	1.47	1.40
4	B	1589	ADP	C5-C4	3.38	1.48	1.40

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1591	ADP	N3-C2-N1	-6.56	123.88	128.89
4	B	1589	ADP	N3-C2-N1	-6.06	124.25	128.89
4	A	1591	ADP	C4-C5-N7	-3.40	106.35	109.48

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
4	B	1589	ADP	C4-C5-N7	-2.71	106.99	109.48
4	B	1589	ADP	C2'-C1'-N9	-2.22	110.91	114.29
4	A	1591	ADP	O3B-PB-O1B	2.11	117.36	110.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 6 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	1590	GOL	3	0
4	B	1589	ADP	3	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, 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	374/392 (95%)	0.02	5 (1%) 79 81	23, 47, 77, 97	0
1	B	374/392 (95%)	0.04	1 (0%) 94 95	31, 52, 87, 132	0
1	C	268/392 (68%)	0.37	18 (6%) 21 22	39, 84, 131, 158	0
1	D	270/392 (68%)	0.96	59 (21%) 1 1	57, 86, 118, 181	0
All	All	1286/1568 (82%)	0.30	83 (6%) 22 24	23, 62, 113, 181	0

All (83) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	430	PHE	5.9
1	D	512	PHE	5.6
1	D	458	LEU	5.5
1	D	515	PRO	5.2
1	D	504	VAL	4.9
1	D	485	ILE	4.9
1	D	513	LEU	4.5
1	D	271	ILE	4.3
1	C	282	ARG	4.1
1	D	524	LEU	4.1
1	C	522	GLN	3.9
1	D	489	ARG	3.8
1	D	257	LEU	3.8
1	D	502	LEU	3.8
1	D	229	TRP	3.7
1	D	240	TRP	3.7
1	D	479	PRO	3.5
1	D	519	ARG	3.4
1	D	283	LYS	3.4
1	D	532	LEU	3.3
1	D	514	ALA	3.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	C	277	TYR	3.3
1	D	277	TYR	3.2
1	D	462	ARG	3.2
1	D	503	ILE	3.1
1	C	438	ALA	3.1
1	D	482	PRO	3.1
1	D	486	ALA	3.0
1	D	300	PRO	3.0
1	D	336	TYR	3.0
1	C	281	CYS	2.9
1	D	523	ALA	2.9
1	D	448	PHE	2.9
1	D	535	ALA	2.9
1	D	236	THR	2.8
1	D	244	ALA	2.8
1	D	525	LEU	2.8
1	D	492	SER	2.8
1	C	427	PHE	2.7
1	C	286	GLY	2.7
1	D	253	THR	2.7
1	D	287	VAL	2.6
1	C	275	ARG	2.6
1	D	455	TRP	2.6
1	D	370	VAL	2.6
1	A	550	LEU	2.6
1	D	457	ILE	2.6
1	D	284	LEU	2.5
1	D	286	GLY	2.5
1	D	475	MET	2.5
1	C	425	SER	2.5
1	C	477	VAL	2.4
1	D	367	PRO	2.4
1	C	498	TYR	2.3
1	D	301	LEU	2.3
1	D	516	GLU	2.3
1	D	429	VAL	2.3
1	A	219	ARG	2.3
1	C	262	TYR	2.3
1	C	520	TYR	2.3
1	D	474	TRP	2.3
1	D	279	GLU	2.3
1	D	290	ASN	2.3

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	D	241	THR	2.3
1	D	533	LEU	2.3
1	C	521	GLY	2.3
1	D	520	TYR	2.2
1	C	495	ILE	2.2
1	C	519	ARG	2.2
1	D	526	ASP	2.2
1	A	561	ALA	2.2
1	B	296	GLY	2.1
1	D	337	LEU	2.1
1	D	453	PHE	2.1
1	D	521	GLY	2.1
1	D	270	ILE	2.1
1	D	522	GLN	2.1
1	A	391	GLU	2.1
1	C	523	ALA	2.0
1	D	231	ARG	2.0
1	A	567	LEU	2.0
1	D	329	LEU	2.0
1	C	428	ILE	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, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
2	MG	C	1533	1/1	0.84	0.25	1.71	77,77,77,77	0

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(\AA^2)	Q<0.9
4	ADP	A	1591	27/27	0.96	0.20	0.97	38,55,73,74	0
4	ADP	B	1589	27/27	0.94	0.16	-0.16	45,59,72,74	0
2	MG	B	1588	1/1	0.86	0.11	-1.81	70,70,70,70	0
3	GOL	A	1593	6/6	0.83	0.20	-	73,79,85,86	0
3	GOL	B	1590	6/6	0.64	0.17	-	90,92,93,93	0
3	GOL	A	1590	6/6	0.81	0.16	-	68,73,76,76	0
2	MG	A	1589	1/1	0.88	0.10	-	50,50,50,50	0
5	SO4	A	1592	5/5	0.91	0.13	-	114,115,117,118	0

6.5 Other polymers

There are no such residues in this entry.