WORLDWIDE PDB PROTEIN DATA BANK

Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 03:40 PM GMT

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 (i)

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

The reported resolution of this entry is 3.50 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.
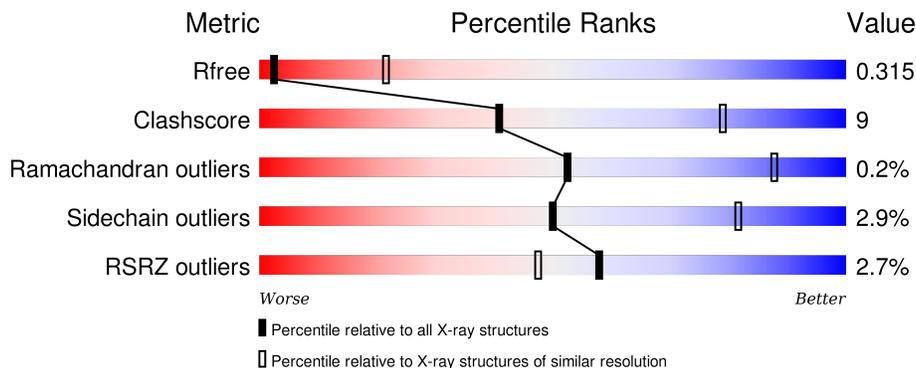
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1051 (3.60-3.40)
Clashscore	102246	1157 (3.60-3.40)
Ramachandran outliers	100387	1120 (3.60-3.40)
Sidechain outliers	100360	1121 (3.60-3.40)
RSRZ outliers	91569	1058 (3.60-3.40)

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	426	2% · 81% · 17% ·
1	B	426	2% · 77% · 21% ··
1	D	426	4% · 74% · 23% ··
2	C	426	3% · 68% · 25% · 5%

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 12548 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 NA(+)/H(+) ANTIPORTER 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	415	Total	C	N	O	S	0	0	0
			3138	2103	473	546	16			
1	B	419	Total	C	N	O	S	0	0	0
			3175	2129	479	551	16			
1	D	417	Total	C	N	O	S	0	0	0
			3159	2118	476	549	16			

- Molecule 2 is a protein called NA(+)/H(+) ANTIPORTER 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	404	Total	C	N	O	S	0	0	0
			3063	2054	461	533	15			

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	367	GLU	ASP	CONFLICT	UNP Q60362

- Molecule 3 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	1	Total	K	0	0
			1	1		
3	C	1	Total	K	0	0
			1	1		

- Molecule 4 is TRIS(HYDROXYETHYL)AMINOMETHANE (three-letter code: TAM) (formula: $C_7H_{17}NO_3$).

TAM
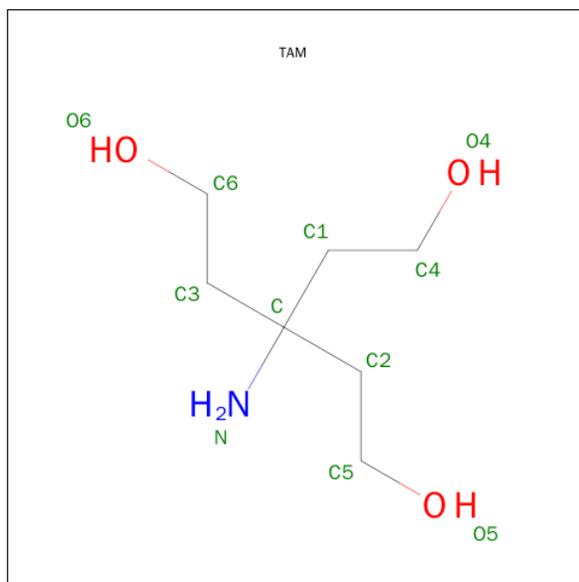
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
4	B	1	Total	C	N	O	0	0
			11	7	1	3		

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: NA(+)/H(+) ANTIPORTER 1
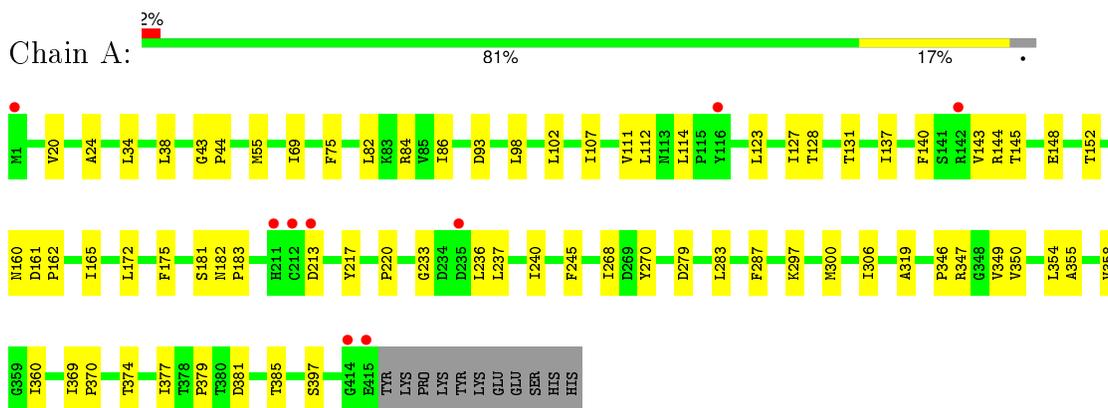
- Molecule 1: NA(+)/H(+) ANTIPORTER 1
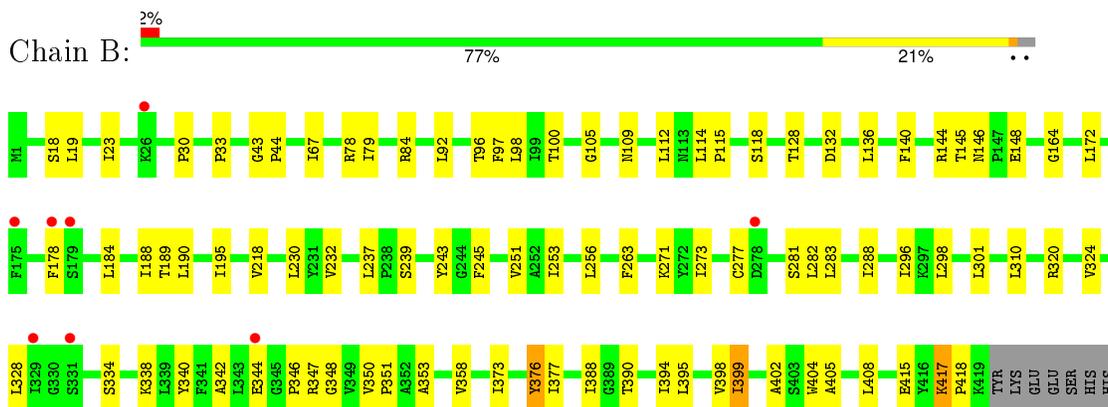
- Molecule 1: NA(+)/H(+) ANTIPORTER 1
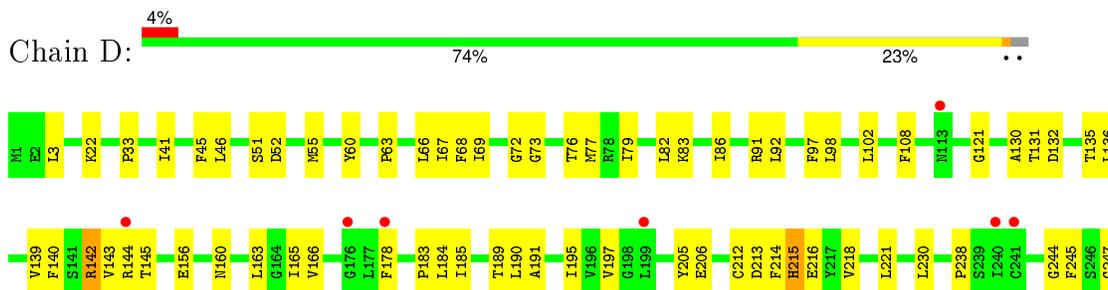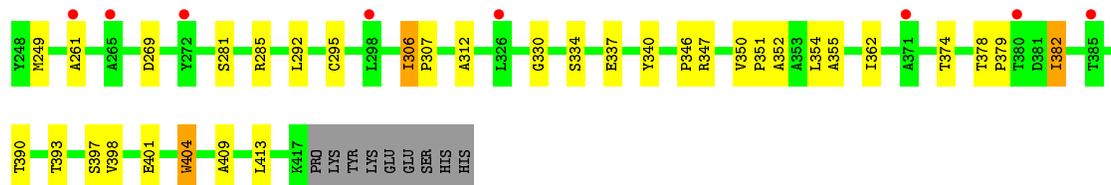
• Molecule 2: NA(+)/H(+) ANTIPORTER 1

Chain C:
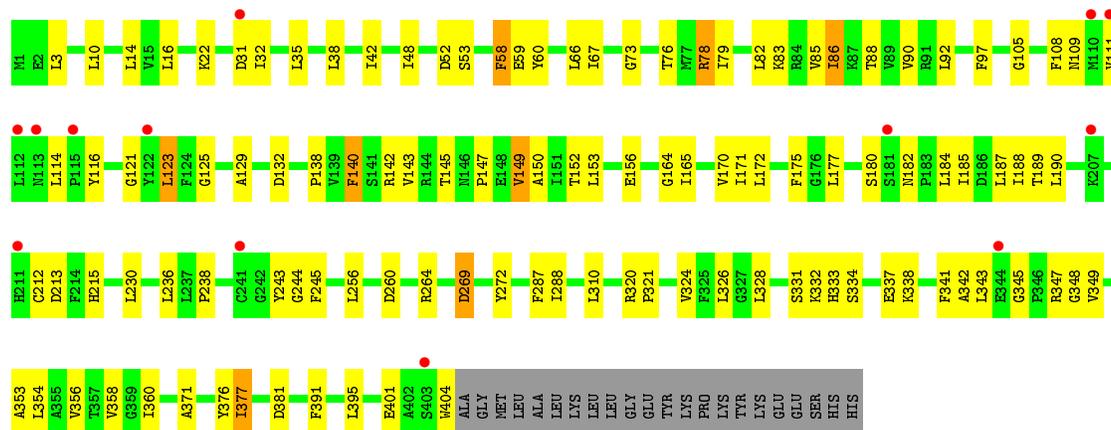
4 Data and refinement statistics (i)

| Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	98.44Å 102.54Å 132.10Å 90.00° 105.63° 90.00°	Depositor
Resolution (Å)	32.00 – 3.50 45.10 – 3.50	Depositor EDS
% Data completeness (in resolution range)	99.3 (32.00-3.50) 97.7 (45.10-3.50)	Depositor EDS
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
$<I/\sigma(I)>$ [1]	1.58 (at 3.48Å)	Xtriage
Refinement program	PHENIX (PHENIX.REFINE)	Depositor
R, R_{free}	0.252 , 0.302 0.268 , 0.315	Depositor DCC
R_{free} test set	1589 reflections (5.33%)	DCC
Wilson B-factor (Å^2)	119.1	Xtriage
Anisotropy	0.283	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.28 , 104.5	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning[2]	$<|L|> = 0.49, <L^2> = 0.32$	Xtriage
Outliers	1 of 31390 reflections (0.003%)	Xtriage
F_o,F_c correlation	0.90	EDS
Total number of atoms	12548	wwPDB-VP
Average B, all atoms (Å^2)	141.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.63% of the height of the origin peak. No significant pseudotranslation is detected.*

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[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 (i)

5.1 Standard geometry (i)

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

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	0.24	0/3202	0.42	0/4357
1	B	0.24	0/3241	0.42	0/4409
1	D	0.24	0/3224	0.43	0/4386
2	C	0.24	0/3127	0.44	0/4257
All	All	0.24	0/12794	0.43	0/17409

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 (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	3138	0	3356	45	0
1	B	3175	0	3398	50	0
1	D	3159	0	3378	66	0
2	C	3063	0	3270	71	0
3	A	1	0	0	0	0
3	C	1	0	0	0	0
4	B	11	0	17	2	0
All	All	12548	0	13419	223	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 9.

All (223) 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:B:190:LEU:HA	1:B:245:PHE:HA	1.66	0.78
1:D:82:LEU:HD13	1:D:86:ILE:HD12	1.70	0.74
1:D:142:ARG:O	1:D:144:ARG:N	2.19	0.73
2:C:78:ARG:HG3	2:C:260:ASP:HB2	1.70	0.71
1:A:69:ILE:HG12	1:A:349:VAL:HG11	1.76	0.68
1:D:306:ILE:HG23	1:D:307:PRO:HD3	1.76	0.68
2:C:59:GLU:HB3	2:C:360:ILE:HD13	1.77	0.67
2:C:172:LEU:HD11	2:C:358:VAL:HG22	1.75	0.67
1:D:214:PHE:HB2	1:D:218:VAL:HG22	1.78	0.66
1:D:108:PHE:HD2	1:D:121:GLY:HA2	1.64	0.63
2:C:238:PRO:HB3	2:C:243:TYR:HA	1.80	0.63
1:D:374:THR:HG23	1:D:379:PRO:HD3	1.81	0.63
1:D:135:THR:HG21	1:D:346:PRO:HG2	1.81	0.62
1:A:165:ILE:HG12	1:A:354:LEU:HD21	1.81	0.62
1:B:288:ILE:HD13	1:B:353:ALA:HB2	1.81	0.62
1:A:268:ILE:HD11	1:D:142:ARG:HB3	1.82	0.62
1:B:105:GLY:O	1:B:109:ASN:ND2	2.33	0.62
1:D:212:CYS:SG	1:D:213:ASP:N	2.74	0.61
2:C:170:VAL:HG21	2:C:187:LEU:HD12	1.83	0.60
1:D:165:ILE:HG12	1:D:354:LEU:HD21	1.84	0.60
1:B:145:THR:OG1	1:B:146:ASN:N	2.34	0.60
2:C:190:LEU:HA	2:C:245:PHE:HA	1.84	0.60
2:C:60:TYR:HD1	1:D:3:LEU:HD13	1.66	0.59
2:C:331:SER:HG	2:C:333:HIS:HD1	1.50	0.59
1:A:112:LEU:HB2	1:A:114:LEU:HD13	1.85	0.59
2:C:328:LEU:HD22	2:C:338:LYS:HG2	1.85	0.59
2:C:345:GLY:O	2:C:347:ARG:NH2	2.36	0.58
2:C:78:ARG:HH21	2:C:256:LEU:HB3	1.69	0.58
1:B:189:THR:HG21	1:B:243:TYR:HB3	1.85	0.58
1:D:409:ALA:HA	1:D:413:LEU:HD13	1.86	0.58
1:A:148:GLU:O	1:A:152:THR:OG1	2.17	0.57
2:C:371:ALA:HA	2:C:377:ILE:HG21	1.86	0.57
1:A:84:ARG:NE	1:A:148:GLU:OE2	2.37	0.57
1:A:24:ALA:HB2	1:A:34:LEU:HD12	1.87	0.57
1:D:68:PHE:O	1:D:72:GLY:N	2.37	0.57
1:B:136:LEU:HD23	1:B:346:PRO:HD2	1.87	0.56
1:D:281:SER:OG	1:D:285:ARG:NH1	2.37	0.56
2:C:165:ILE:HD13	2:C:354:LEU:HD11	1.88	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:16:LEU:HB3	2:C:38:LEU:HD21	1.86	0.56
1:B:277:CYS:O	1:B:281:SER:N	2.32	0.56
2:C:321:PRO:HA	2:C:342:ALA:HB1	1.88	0.55
2:C:92:LEU:HD22	2:C:156:GLU:HG3	1.88	0.55
1:B:67:ILE:HD11	1:B:251:VAL:HG21	1.88	0.55
2:C:132:ASP:HB2	2:C:348:GLY:HA3	1.89	0.55
2:C:125:GLY:O	2:C:129:ALA:N	2.40	0.54
2:C:67:ILE:HD11	2:C:230:LEU:HD13	1.89	0.54
1:D:362:ILE:HD13	1:D:382:ILE:HD11	1.90	0.54
2:C:52:ASP:OD1	2:C:53:SER:N	2.40	0.54
2:C:269:ASP:OD1	2:C:269:ASP:N	2.40	0.54
1:D:401:GLU:HA	1:D:404:TRP:HB2	1.88	0.54
2:C:111:VAL:HG23	2:C:310:LEU:HD23	1.89	0.54
1:D:41:ILE:HA	1:D:45:PHE:HB2	1.89	0.54
1:A:143:VAL:HG23	1:A:145:THR:HG23	1.89	0.53
1:D:136:LEU:HD22	1:D:139:VAL:HB	1.90	0.53
1:D:404:TRP:HA	1:D:404:TRP:CE3	2.43	0.53
1:D:205:TYR:OH	1:D:221:LEU:O	2.20	0.53
1:B:218:VAL:HG11	1:B:273:ILE:HG12	1.92	0.52
2:C:22:LYS:HD3	1:D:214:PHE:CE1	2.44	0.52
1:D:183:PRO:HG2	1:D:184:LEU:HD12	1.91	0.52
1:B:84:ARG:HE	1:B:148:GLU:HG3	1.74	0.52
1:B:334:SER:OG	1:B:415:GLU:OE2	2.18	0.52
1:D:131:THR:OG1	1:D:160:ASN:OD1	2.25	0.52
2:C:180:SER:O	2:C:182:ASN:ND2	2.43	0.52
1:D:218:VAL:HG21	1:D:269:ASP:HB3	1.90	0.52
1:D:404:TRP:HE3	1:D:404:TRP:HA	1.75	0.51
1:D:91:ARG:HH22	1:D:330:GLY:H	1.59	0.51
1:D:92:LEU:HD23	1:D:156:GLU:HG3	1.91	0.51
1:D:136:LEU:O	1:D:140:PHE:N	2.42	0.51
1:D:97:PHE:HD2	1:D:98:LEU:HD12	1.76	0.51
1:D:33:PRO:HD3	1:D:398:VAL:HG11	1.93	0.51
1:B:232:VAL:O	1:B:237:LEU:HB2	2.11	0.51
2:C:123:LEU:HB3	2:C:171:ILE:HG21	1.93	0.50
1:B:328:LEU:O	1:B:338:LYS:NZ	2.43	0.50
2:C:31:ASP:N	2:C:31:ASP:OD1	2.43	0.50
1:B:390:THR:O	1:B:394:ILE:HG12	2.12	0.50
2:C:85:VAL:HG22	2:C:88:THR:HB	1.94	0.50
1:A:127:ILE:HD13	1:A:355:ALA:HA	1.94	0.50
2:C:78:ARG:HD2	2:C:79:ILE:H	1.77	0.49
1:A:220:PRO:HB2	1:B:18:SER:HB3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:417:LYS:HE2	2:C:142:ARG:HD3	1.94	0.49
2:C:86:ILE:O	2:C:90:VAL:HG22	2.12	0.49
1:A:131:THR:OG1	1:A:160:ASN:OD1	2.29	0.49
1:D:238:PRO:HD3	1:D:244:GLY:HA3	1.94	0.48
1:A:93:ASP:OD1	1:A:160:ASN:ND2	2.46	0.48
1:B:417:LYS:HD3	1:B:417:LYS:H	1.78	0.48
2:C:109:ASN:HA	2:C:114:LEU:HB2	1.95	0.48
1:D:51:SER:OG	1:D:295:CYS:O	2.29	0.48
1:D:230:LEU:HA	1:D:247:GLY:HA3	1.94	0.48
1:B:376:TYR:CE2	1:B:377:ILE:HG12	2.49	0.48
2:C:78:ARG:HD2	2:C:79:ILE:HG23	1.94	0.48
2:C:149:VAL:HG23	2:C:341:PHE:HE1	1.79	0.48
2:C:331:SER:OG	2:C:333:HIS:ND1	2.39	0.47
1:D:334:SER:HB3	1:D:337:GLU:HG3	1.95	0.47
1:D:166:VAL:HG13	1:D:190:LEU:HD12	1.95	0.47
2:C:58:PHE:CD2	2:C:356:VAL:HG11	2.50	0.47
1:B:373:ILE:H	1:B:373:ILE:HD12	1.79	0.47
1:A:137:ILE:HA	1:A:140:PHE:CD1	2.49	0.47
1:B:128:THR:O	1:B:347:ARG:NE	2.43	0.47
1:B:395:LEU:O	1:B:399:ILE:HG22	2.13	0.47
1:B:79:ILE:HG21	1:B:256:LEU:HD21	1.97	0.47
1:B:324:VAL:HG21	1:B:342:ALA:HA	1.97	0.47
1:B:418:PRO:HG2	2:C:138:PRO:HB3	1.97	0.47
1:D:66:LEU:HA	1:D:69:ILE:HD12	1.97	0.47
1:B:19:LEU:O	1:B:23:ILE:HG12	2.15	0.47
2:C:260:ASP:OD2	2:C:264:ARG:NH1	2.49	0.46
2:C:238:PRO:HD3	2:C:244:GLY:HA3	1.97	0.46
1:D:130:ALA:HA	1:D:160:ASN:HB3	1.97	0.46
2:C:3:LEU:HD22	1:D:60:TYR:HA	1.96	0.46
2:C:140:PHE:HA	2:C:143:VAL:HG12	1.98	0.46
1:A:137:ILE:HA	1:A:140:PHE:HD1	1.79	0.46
1:D:215:HIS:HB3	1:D:216:GLU:H	1.62	0.46
1:D:340:TYR:CG	1:D:413:LEU:HD11	2.51	0.46
1:D:132:ASP:HB3	1:D:135:THR:HG23	1.96	0.46
1:D:41:ILE:HG23	1:D:46:LEU:HD13	1.98	0.46
2:C:401:GLU:HA	2:C:404:TRP:CD1	2.51	0.46
1:A:374:THR:HG23	1:A:379:PRO:HD3	1.97	0.45
4:B:1420:TAM:H42	4:B:1420:TAM:H21	1.52	0.45
1:D:190:LEU:HA	1:D:245:PHE:HA	1.98	0.45
1:A:233:GLY:HA2	1:A:245:PHE:HB3	1.99	0.45
1:B:140:PHE:O	1:B:144:ARG:N	2.45	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:172:LEU:HB3	2:C:177:LEU:HD12	1.99	0.45
2:C:401:GLU:HA	2:C:404:TRP:HD1	1.80	0.45
2:C:147:PRO:HA	2:C:150:ALA:HB3	1.99	0.45
1:A:112:LEU:HD11	1:A:385:THR:HG21	1.99	0.45
1:B:132:ASP:HB2	1:B:348:GLY:HA3	1.99	0.45
1:A:279:ASP:O	1:A:283:LEU:HG	2.17	0.44
2:C:42:ILE:HA	2:C:48:ILE:HD12	1.99	0.44
1:A:346:PRO:HA	1:A:397:SER:OG	2.17	0.44
1:A:161:ASP:HB3	1:A:350:VAL:HG21	1.98	0.44
2:C:32:ILE:HA	2:C:35:LEU:HD23	2.00	0.44
1:A:306:ILE:HD12	1:A:306:ILE:H	1.83	0.44
1:B:30:PRO:HG2	1:B:402:ALA:HB2	1.99	0.44
2:C:79:ILE:HA	2:C:82:LEU:HB2	1.99	0.44
2:C:288:ILE:HG21	2:C:349:VAL:HG23	1.99	0.44
1:D:292:LEU:HD22	1:D:352:ALA:HB1	1.98	0.44
1:A:236:LEU:O	1:A:240:ILE:HG12	2.18	0.44
1:D:378:THR:O	1:D:382:ILE:HG23	2.18	0.44
1:B:195:ILE:HD13	1:B:253:ILE:HG13	1.99	0.44
1:B:115:PRO:HG2	1:B:118:SER:HB2	2.00	0.44
1:B:100:THR:HG22	1:B:320:ARG:HA	2.00	0.44
1:A:370:PRO:O	1:A:374:THR:OG1	2.28	0.44
2:C:108:PHE:HD2	2:C:121:GLY:HA2	1.83	0.43
1:A:75:PHE:O	1:A:270:TYR:OH	2.32	0.43
1:A:220:PRO:HG3	1:B:283:LEU:HD23	1.99	0.43
1:D:197:VAL:HG11	1:D:245:PHE:CZ	2.53	0.43
2:C:66:LEU:HD11	2:C:353:ALA:HB1	2.01	0.43
2:C:334:SER:HB3	2:C:337:GLU:HG2	1.99	0.43
1:A:369:ILE:HD12	1:A:370:PRO:HD2	2.00	0.43
2:C:391:PHE:CE1	2:C:395:LEU:HD11	2.54	0.43
2:C:324:VAL:HG13	2:C:328:LEU:HD13	2.01	0.43
1:D:98:LEU:O	1:D:102:LEU:HG	2.19	0.43
1:A:297:LYS:HB2	1:A:300:MET:HG2	2.01	0.43
1:A:172:LEU:HA	1:A:172:LEU:HD23	1.92	0.43
1:A:128:THR:HG23	1:A:347:ARG:HG2	2.00	0.43
1:A:114:LEU:HD11	1:A:377:ILE:HG13	2.01	0.43
1:B:97:PHE:HD2	1:B:98:LEU:HD12	1.84	0.43
1:B:376:TYR:CZ	1:B:377:ILE:HG12	2.54	0.43
1:B:301:LEU:HA	1:B:388:ILE:HD13	2.00	0.43
2:C:73:GLY:O	2:C:76:THR:OG1	2.27	0.43
1:D:163:LEU:HD23	1:D:249:MET:HE1	2.00	0.42
2:C:78:ARG:HA	2:C:78:ARG:HD3	1.91	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:381:ASP:O	1:A:385:THR:OG1	2.31	0.42
1:A:213:ASP:N	1:A:213:ASP:OD1	2.44	0.42
1:D:312:ALA:HB1	1:D:393:THR:HA	2.01	0.42
1:A:20:VAL:HG21	1:A:38:LEU:HD22	2.01	0.42
1:D:82:LEU:HD12	1:D:83:LYS:N	2.35	0.42
2:C:60:TYR:HA	1:D:3:LEU:HD22	2.00	0.42
2:C:10:LEU:O	2:C:14:LEU:HG	2.20	0.42
1:D:185:ILE:O	1:D:189:THR:HG23	2.19	0.42
1:B:67:ILE:HD13	1:B:230:LEU:HD11	2.00	0.42
1:D:41:ILE:O	1:D:46:LEU:HB2	2.20	0.42
1:A:43:GLY:HA3	1:A:44:PRO:HD3	1.92	0.42
1:D:63:PRO:O	1:D:67:ILE:HG13	2.19	0.42
1:D:346:PRO:HA	1:D:397:SER:HB2	2.02	0.42
2:C:85:VAL:HG11	2:C:152:THR:HG22	2.02	0.42
1:A:82:LEU:O	1:A:86:ILE:N	2.53	0.42
1:D:347:ARG:HB3	1:D:351:PRO:HG2	2.00	0.42
1:A:98:LEU:O	1:A:102:LEU:HG	2.20	0.42
2:C:185:ILE:O	2:C:189:THR:HG23	2.20	0.42
1:B:172:LEU:HD11	1:B:358:VAL:HG22	2.00	0.42
1:D:69:ILE:O	1:D:73:GLY:N	2.49	0.42
2:C:105:GLY:HA2	2:C:121:GLY:C	2.40	0.41
1:B:112:LEU:HB2	1:B:114:LEU:HD13	2.02	0.41
1:A:55:MET:HB3	1:A:360:ILE:HG22	2.01	0.41
1:A:137:ILE:HG13	1:A:140:PHE:HE1	1.85	0.41
1:B:44:PRO:HB3	1:B:298:LEU:HB2	2.01	0.41
1:B:43:GLY:HA3	1:B:44:PRO:HD3	1.88	0.41
1:D:206:GLU:HG3	1:D:261:ALA:HB1	2.01	0.41
1:D:205:TYR:CE2	1:D:221:LEU:HG	2.55	0.41
2:C:184:LEU:O	2:C:188:ILE:HG12	2.21	0.41
1:D:350:VAL:HB	1:D:351:PRO:HD3	2.02	0.41
1:B:44:PRO:HG2	1:B:296:ILE:HG22	2.03	0.41
1:D:52:ASP:O	1:D:55:MET:HB2	2.21	0.41
1:D:83:LYS:HA	1:D:83:LYS:HD3	1.88	0.41
1:D:108:PHE:CD2	1:D:121:GLY:HA2	2.51	0.41
2:C:343:LEU:HB3	2:C:404:TRP:CZ2	2.56	0.41
1:A:107:ILE:O	1:A:111:VAL:HG23	2.21	0.41
1:B:92:LEU:O	1:B:96:THR:HB	2.20	0.41
1:B:340:TYR:O	1:B:344:GLU:HB2	2.19	0.41
1:B:402:ALA:O	4:B:1420:TAM:N	2.35	0.41
1:B:97:PHE:HZ	1:B:164:GLY:HA2	1.85	0.41
2:C:212:CYS:SG	2:C:213:ASP:N	2.93	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:181:SER:OG	1:A:182:ASN:N	2.54	0.41
2:C:172:LEU:HA	2:C:175:PHE:HB3	2.03	0.41
2:C:79:ILE:O	2:C:83:LYS:HG2	2.21	0.40
2:C:332:LYS:HD2	2:C:332:LYS:HA	1.86	0.40
1:B:33:PRO:HD3	1:B:398:VAL:HG11	2.03	0.40
1:B:417:LYS:CD	1:B:417:LYS:H	2.32	0.40
1:D:355:ALA:HB2	1:D:390:THR:HG21	2.02	0.40
1:A:182:ASN:HA	1:A:183:PRO:HD3	1.93	0.40
2:C:97:PHE:HZ	2:C:164:GLY:HA2	1.85	0.40
2:C:320:ARG:HB3	2:C:321:PRO:HD3	2.04	0.40
2:C:150:ALA:HA	2:C:153:LEU:HD13	2.03	0.40
1:A:161:ASP:HB2	1:A:162:PRO:HD3	2.03	0.40
2:C:108:PHE:CD2	2:C:121:GLY:HA2	2.57	0.40
1:B:184:LEU:O	1:B:188:ILE:HG12	2.21	0.40
2:C:215:HIS:HB2	2:C:272:TYR:CZ	2.56	0.40
1:D:191:ALA:O	1:D:195:ILE:HG12	2.21	0.40
1:A:123:LEU:HD21	1:A:358:VAL:HG11	2.04	0.40
1:A:175:PHE:CD2	1:A:369:ILE:HD11	2.56	0.40
1:B:350:VAL:HB	1:B:351:PRO:HD3	2.03	0.40
1:A:217:TYR:CD1	1:B:282:LEU:HD21	2.57	0.40
1:D:22:LYS:HA	1:D:22:LYS:HD2	1.87	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	413/426 (97%)	385 (93%)	27 (6%)	1 (0%)	52	88
1	B	417/426 (98%)	383 (92%)	33 (8%)	1 (0%)	52	88
1	D	415/426 (97%)	384 (92%)	30 (7%)	1 (0%)	52	88
2	C	402/426 (94%)	366 (91%)	36 (9%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1647/1704 (97%)	1518 (92%)	126 (8%)	3 (0%)	52 88

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	143	VAL
1	B	405	ALA
1	A	319	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	340/351 (97%)	337 (99%)	3 (1%)	84 94
1	B	344/351 (98%)	333 (97%)	11 (3%)	46 80
1	D	342/351 (97%)	332 (97%)	10 (3%)	50 81
2	C	333/351 (95%)	318 (96%)	15 (4%)	34 73
All	All	1359/1404 (97%)	1320 (97%)	39 (3%)	50 81

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

Mol	Chain	Res	Type
1	A	144	ARG
1	A	237	LEU
1	A	287	PHE
1	B	78	ARG
1	B	178	PHE
1	B	239	SER
1	B	263	PHE
1	B	271	LYS
1	B	310	LEU
1	B	376	TYR
1	B	399	ILE
1	B	404	TRP

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Mol	Chain	Res	Type
1	B	408	LEU
1	B	417	LYS
2	C	58	PHE
2	C	78	ARG
2	C	86	ILE
2	C	116	TYR
2	C	123	LEU
2	C	140	PHE
2	C	145	THR
2	C	149	VAL
2	C	236	LEU
2	C	269	ASP
2	C	287	PHE
2	C	326	LEU
2	C	376	TYR
2	C	377	ILE
2	C	381	ASP
1	D	76	THR
1	D	77	MET
1	D	79	ILE
1	D	142	ARG
1	D	145	THR
1	D	178	PHE
1	D	215	HIS
1	D	306	ILE
1	D	382	ILE
1	D	404	TRP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 3 ligands modelled in this entry, 2 are monoatomic - leaving 1 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	$\#|Z| > 2$
4	TAM	B	1420	-	7,10,10	1.13	0	9,12,12	0.68	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
4	TAM	B	1420	-	-	0/12/12/12	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.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	1420	TAM	2	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 (i)

6.1 Protein, DNA and RNA chains (i)

In the following table, the column labelled '#RSRZ> 2' contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled 'Q< 0.9' lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	415/426 (97%)	-0.21	9 (2%)	65	55	83, 116, 155, 199	0
1	B	419/426 (98%)	-0.11	8 (1%)	70	60	87, 133, 179, 201	0
1	D	417/426 (97%)	0.02	15 (3%)	46	37	105, 145, 186, 210	0
2	C	404/426 (94%)	-0.08	13 (3%)	51	42	106, 162, 192, 220	0
All	All	1655/1704 (97%)	-0.09	45 (2%)	58	47	83, 139, 186, 220	0

All (45) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	240	ILE	4.0
1	D	241	CYS	3.7
1	D	371	ALA	3.6
2	C	110	MET	3.6
1	B	278	ASP	3.5
1	A	235	ASP	3.3
1	A	142	ARG	3.3
2	C	403	SER	3.2
1	D	380	THR	3.1
2	C	111	VAL	3.1
1	A	415	GLU	3.0
1	D	261	ALA	3.0
1	B	329	ILE	3.0
1	D	265	ALA	2.9
2	C	181	SER	2.9
1	B	344	GLU	2.9
2	C	344	GLU	2.9
2	C	113	ASN	2.8
1	A	211	HIS	2.8
1	A	116	TYR	2.7
1	A	1	MET	2.7

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Mol	Chain	Res	Type	RSRZ
1	B	179	SER	2.6
2	C	31	ASP	2.6
1	D	178	PHE	2.6
1	D	326	LEU	2.5
1	A	213	ASP	2.5
1	B	175	PHE	2.4
1	D	176	GLY	2.4
2	C	211	HIS	2.4
1	D	298	LEU	2.3
2	C	207	LYS	2.3
2	C	115	PRO	2.3
1	B	178	PHE	2.3
1	D	385	THR	2.3
1	A	414	GLY	2.3
1	B	331	SER	2.3
1	A	212	CYS	2.2
2	C	122	TYR	2.2
2	C	241	CYS	2.1
2	C	112	LEU	2.1
1	D	199	LEU	2.1
1	D	113	ASN	2.1
1	D	144	ARG	2.1
1	D	272	TYR	2.0
1	B	26	LYS	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
4	TAM	B	1420	11/11	0.77	0.19	-0.93	107,128,151,154	0
3	K	C	1405	1/1	0.91	0.12	-1.36	127,127,127,127	0
3	K	A	1416	1/1	0.86	0.17	-1.66	128,128,128,128	0

6.5 Other polymers (i)

There are no such residues in this entry.