Full wwPDB X-ray Structure Validation Report (i)

Oct 12, 2021 – 01:26 PM EDT

PDB ID	:	2A8M
Title	:	Crystal Structure of Human Taspase1 (T234S mutant)
Authors	:	Khan, J.A.; Dunn, B.M.; Tong, L.
Deposited on	:	2005-07-08
Resolution	:	2.60 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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

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

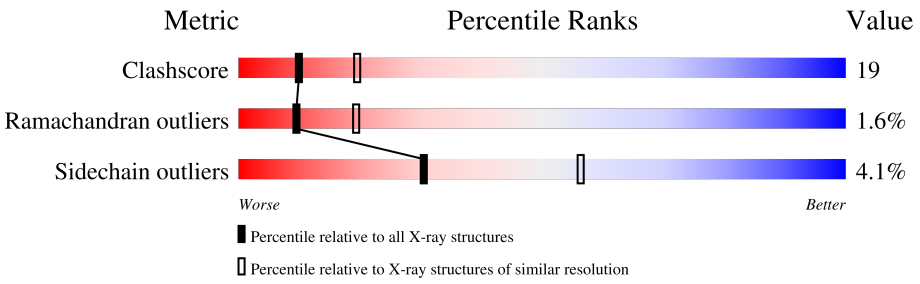
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.23.2

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 2.60 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	141614	3518 (2.60-2.60)
Ramachandran outliers	138981	3455 (2.60-2.60)
Sidechain outliers	138945	3455 (2.60-2.60)

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 of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. 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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	420	51% 22% • 26%
1	B	420	45% 28% • 25%

2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4623 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 Threonine aspartase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	310	2246	1399	400	429	18	0	0	0
1	B	313	2266	1409	404	434	19	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	234	SER	THR	engineered mutation	UNP Q9H6P5
B	234	SER	THR	engineered mutation	UNP Q9H6P5

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
3	A	53	53	53	0	0
3	B	56	56	56	0	0

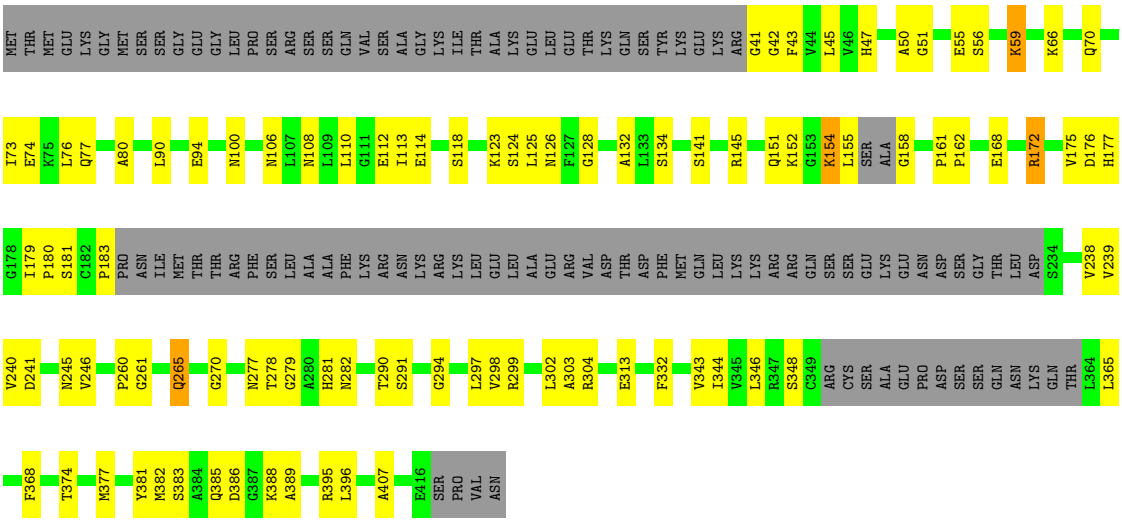
3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. 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. 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.

Note EDS was not executed.

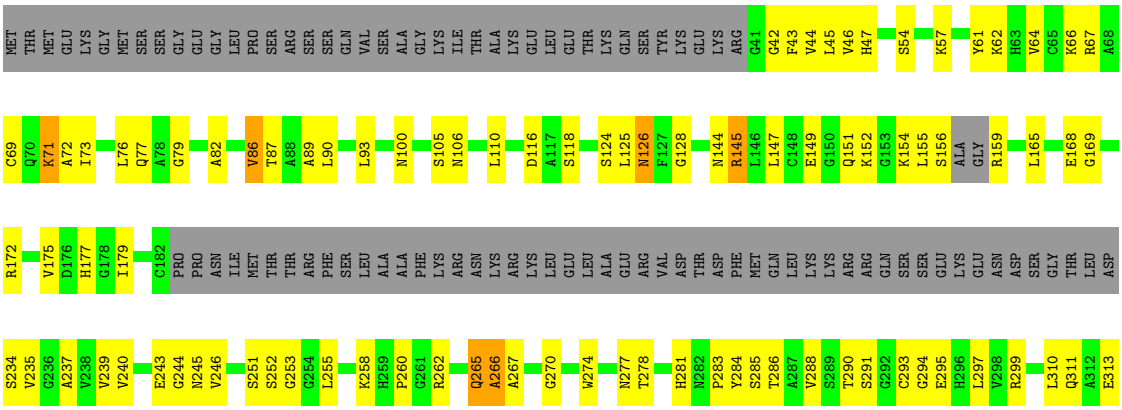
• Molecule 1: Threonine aspartase 1

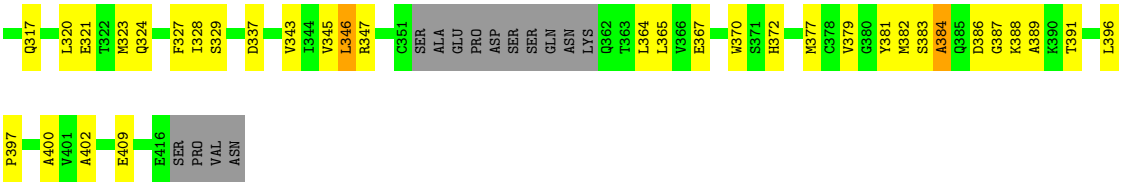
Chain A: 51% 22% • 26%



• Molecule 1: Threonine aspartase 1

Chain B: 45% 28% • 25%




Q317
L320
E321
T322
M323
Q324
F327
I328
S329
D337
V343
I344
V345
L346
R347
C351
SER
ALA
GLU
PRO
ASP
SER
SER
GLN
ASN
LYS
Q362
T363
L364
L365
V366
E367
W370
S371
H372
M377
C378
V379
G380
Y381
M382
S383
A384
Q385
D386
G387
K388
A389
K390
T391
L396
P397
A400
V401
A402
E409
E416
SER
PRO
VAL
ASN

4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	60.44Å 89.37Å 104.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.59 – 2.60	Depositor
% Data completeness (in resolution range)	88.4 (29.59-2.60)	Depositor
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.207 , 0.290	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	4623	wwPDB-VP
Average B, all atoms (Å^2)	45.0	wwPDB-VP

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: CL

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	0.38	0/2284	0.60	0/3088
1	B	0.36	0/2303	0.59	0/3113
All	All	0.37	0/4587	0.59	0/6201

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	2246	0	2219	75	0
1	B	2266	0	2232	103	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	53	0	0	0	0
3	B	56	0	0	5	0
All	All	4623	0	4451	171	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 19.

All (171) 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:383:SER:HB3	1:A:386:ASP:HB2	1.51	0.90
1:A:90:LEU:O	1:A:94:GLU:HG3	1.78	0.81
1:A:260:PRO:O	1:B:169:GLY:HA3	1.80	0.81
1:A:100:ASN:ND2	1:A:106:ASN:H	1.80	0.80
1:B:145:ARG:HG3	1:B:179:ILE:HD11	1.64	0.79
1:A:126:ASN:HD22	1:A:151:GLN:HG2	1.47	0.78
1:A:110:LEU:HD11	1:B:172:ARG:NH1	2.00	0.76
1:B:116:ASP:HB3	1:B:267:ALA:HB2	1.68	0.74
1:A:396:LEU:HD23	1:A:407:ALA:HB2	1.70	0.72
1:B:42:GLY:H	1:B:384:ALA:HB2	1.54	0.72
1:A:299:ARG:HH11	1:A:299:ARG:HG3	1.56	0.70
1:B:383:SER:HB3	1:B:386:ASP:HB2	1.74	0.69
1:A:145:ARG:NH1	1:A:177:HIS:HB3	2.07	0.69
1:B:100:ASN:ND2	1:B:106:ASN:H	1.92	0.67
1:A:290:THR:HG22	1:A:343:VAL:HG12	1.75	0.67
1:A:70:GLN:HA	1:A:382:MET:HE3	1.77	0.66
1:A:100:ASN:HD21	1:A:106:ASN:H	1.43	0.65
1:B:265:GLN:HG2	1:B:294:GLY:HA2	1.78	0.64
1:A:45:LEU:HB3	1:A:238:VAL:CG2	2.27	0.64
1:B:288:VAL:HG13	1:B:345:VAL:HG12	1.78	0.64
1:B:87:THR:OG1	1:B:144:ASN:ND2	2.31	0.64
1:B:365:LEU:HD12	3:B:433:HOH:O	1.98	0.62
1:A:59:LYS:HA	1:A:59:LYS:HE3	1.81	0.62
1:B:42:GLY:HA3	1:B:76:LEU:HD21	1.82	0.62
1:B:370:TRP:NE1	1:B:409:GLU:HB2	2.14	0.61
1:B:62:LYS:O	1:B:66:LYS:HG3	2.00	0.60
1:A:332:PHE:CE2	1:B:125:LEU:HD22	2.36	0.60
1:A:41:GLY:HA2	1:A:385:GLN:NE2	2.18	0.59
1:A:73:ILE:O	1:A:77:GLN:HG3	2.02	0.59
1:A:41:GLY:HA2	1:A:385:GLN:HE21	1.68	0.58
1:A:108:ASN:HD21	1:A:112:GLU:HB2	1.68	0.58
1:B:310:LEU:HD11	1:B:345:VAL:HG21	1.84	0.58
1:B:290:THR:HB	1:B:297:LEU:HD22	1.85	0.58
1:A:42:GLY:HA3	1:A:76:LEU:HD13	1.86	0.57
1:A:70:GLN:O	1:A:74:GLU:HG3	2.05	0.56
1:B:290:THR:HG22	1:B:343:VAL:HG12	1.87	0.56
1:B:62:LYS:HE2	1:B:391:THR:OG1	2.06	0.56
1:A:55:GLU:HA	1:A:55:GLU:OE1	2.06	0.56
1:B:246:VAL:HG22	1:B:277:ASN:HD22	1.71	0.56
1:B:370:TRP:CE3	1:B:379:VAL:HG21	2.41	0.56
1:B:73:ILE:HG13	1:B:382:MET:SD	2.46	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:155:LEU:HD12	1:B:159:ARG:HG2	1.88	0.55
1:B:346:LEU:C	1:B:346:LEU:HD12	2.27	0.55
1:B:382:MET:CE	1:B:389:ALA:HB2	2.36	0.55
1:A:47:HIS:HB3	1:A:344:ILE:HG21	1.88	0.55
1:A:168:GLU:O	1:A:172:ARG:HG3	2.06	0.55
1:B:105:SER:HB3	1:B:253:GLY:HA3	1.88	0.55
1:A:348:SER:HA	1:A:365:LEU:O	2.06	0.55
1:B:239:VAL:CG1	1:B:240:VAL:N	2.70	0.55
1:A:382:MET:HB2	1:A:389:ALA:HA	1.89	0.55
1:B:77:GLN:C	1:B:79:GLY:H	2.10	0.54
1:B:47:HIS:CE1	1:B:291:SER:HB2	2.42	0.54
1:B:370:TRP:CE2	1:B:409:GLU:HB2	2.43	0.54
1:A:299:ARG:HG3	1:A:299:ARG:NH1	2.22	0.54
1:B:155:LEU:HB2	1:B:159:ARG:HG2	1.88	0.54
1:B:86:VAL:O	1:B:90:LEU:HD23	2.08	0.54
1:A:45:LEU:HD13	1:A:381:TYR:HB3	1.90	0.53
1:B:67:ARG:O	1:B:71:LYS:HD2	2.08	0.53
1:B:69:CYS:O	1:B:72:ALA:HB3	2.09	0.53
1:B:246:VAL:HG22	1:B:277:ASN:ND2	2.24	0.53
1:B:258:LYS:HB2	1:B:262:ARG:HD3	1.90	0.53
1:B:46:VAL:HG12	1:B:237:ALA:CB	2.39	0.53
1:A:152:LYS:O	1:A:155:LEU:HD13	2.08	0.53
1:B:42:GLY:N	1:B:384:ALA:HB2	2.23	0.53
1:A:294:GLY:O	1:A:298:VAL:HG23	2.09	0.53
1:B:45:LEU:HD13	1:B:381:TYR:HB3	1.91	0.53
1:B:320:LEU:O	1:B:324:GLN:HG3	2.09	0.53
1:A:45:LEU:CD1	1:A:381:TYR:HB3	2.40	0.52
1:B:145:ARG:NH2	1:B:149:GLU:OE2	2.43	0.52
1:B:72:ALA:O	1:B:76:LEU:HB2	2.10	0.52
1:B:110:LEU:HD22	1:B:260:PRO:HG2	1.91	0.52
1:A:47:HIS:CE1	1:A:291:SER:HB2	2.45	0.52
1:B:270:GLY:HA2	1:B:274:TRP:CE3	2.45	0.52
1:B:154:LYS:HB3	3:B:441:HOH:O	2.10	0.51
1:B:372:HIS:CD2	1:B:377:MET:HB3	2.46	0.51
1:B:118:SER:HA	1:B:128:GLY:O	2.11	0.51
1:B:386:ASP:OD1	1:B:388:LYS:HG2	2.11	0.50
1:A:66:LYS:HB3	1:A:66:LYS:HZ3	1.76	0.50
1:A:126:ASN:HA	1:B:299:ARG:HH12	1.75	0.50
1:B:77:GLN:C	1:B:79:GLY:N	2.64	0.50
1:A:123:LYS:HE3	1:A:245:ASN:ND2	2.26	0.50
1:A:152:LYS:O	1:A:155:LEU:HD22	2.11	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:62:LYS:O	1:B:391:THR:HG21	2.12	0.50
1:A:145:ARG:HH12	1:A:177:HIS:HB3	1.75	0.50
1:B:154:LYS:HD3	3:B:441:HOH:O	2.10	0.50
1:B:255:LEU:HD23	1:B:262:ARG:NH1	2.26	0.49
1:B:246:VAL:H	1:B:277:ASN:HD21	1.58	0.49
1:B:66:LYS:HB3	1:B:66:LYS:NZ	2.28	0.49
1:A:241:ASP:OD2	1:A:245:ASN:HB2	2.13	0.49
1:B:310:LEU:O	1:B:347:ARG:NH1	2.36	0.49
1:B:90:LEU:CD2	1:B:237:ALA:HB3	2.43	0.48
1:B:234:SER:HB3	1:B:252:SER:HB2	1.94	0.48
1:A:51:GLY:O	1:A:374:THR:HB	2.12	0.48
1:B:323:MET:O	1:B:328:ILE:HG13	2.14	0.48
1:B:278:THR:HG23	1:B:285:SER:OG	2.13	0.48
1:A:113:ILE:O	1:A:134:SER:HB3	2.14	0.48
1:A:270:GLY:O	1:A:303:ALA:HB3	2.14	0.48
1:B:67:ARG:NH1	3:B:449:HOH:O	2.44	0.48
1:A:118:SER:HA	1:A:128:GLY:O	2.14	0.48
1:A:175:VAL:C	1:A:177:HIS:H	2.17	0.48
1:B:90:LEU:CD1	1:B:235:VAL:HG23	2.44	0.47
1:B:317:GLN:O	1:B:321:GLU:HG3	2.14	0.47
1:B:278:THR:HG22	1:B:285:SER:N	2.29	0.47
1:A:110:LEU:HD11	1:B:172:ARG:HH12	1.77	0.47
1:A:279:GLY:C	1:A:281:HIS:H	2.17	0.47
1:B:239:VAL:HG12	1:B:240:VAL:N	2.29	0.47
1:B:244:GLY:HA2	1:B:364:LEU:HD21	1.95	0.47
1:B:286:THR:HG21	1:B:311:GLN:NE2	2.30	0.47
1:A:126:ASN:ND2	1:A:151:GLN:HG2	2.24	0.46
1:A:161:PRO:HA	1:A:162:PRO:HD3	1.77	0.46
1:A:50:ALA:HB3	1:A:291:SER:OG	2.15	0.46
1:A:239:VAL:HG12	1:A:240:VAL:N	2.31	0.46
1:B:124:SER:O	1:B:125:LEU:HB2	2.15	0.46
1:A:278:THR:HG23	1:A:282:ASN:O	2.16	0.46
1:B:54:SER:OG	1:B:57:LYS:HB3	2.15	0.46
1:B:71:LYS:HE3	1:B:71:LYS:HA	1.98	0.46
1:B:126:ASN:HB3	1:B:147:LEU:HG	1.99	0.45
1:B:283:PRO:HD2	1:B:311:GLN:OE1	2.16	0.45
1:A:110:LEU:HD21	1:B:172:ARG:HH12	1.82	0.45
1:A:154:LYS:O	1:A:158:GLY:HA2	2.15	0.45
1:A:66:LYS:HB3	1:A:66:LYS:NZ	2.31	0.45
1:A:168:GLU:OE2	1:A:172:ARG:NE	2.49	0.45
1:B:61:TYR:O	1:B:64:VAL:HG12	2.17	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:346:LEU:C	1:A:346:LEU:HD23	2.38	0.45
1:B:71:LYS:HE3	1:B:71:LYS:CA	2.47	0.45
1:B:281:HIS:O	1:B:283:PRO:HD3	2.16	0.45
1:B:382:MET:HE1	1:B:389:ALA:HB2	1.99	0.44
1:A:45:LEU:HB3	1:A:238:VAL:HG22	1.97	0.44
1:B:66:LYS:HB3	1:B:66:LYS:HZ3	1.81	0.44
1:B:89:ALA:O	1:B:93:LEU:HD22	2.17	0.44
1:B:386:ASP:C	1:B:388:LYS:H	2.21	0.44
1:B:265:GLN:HG2	1:B:297:LEU:HD12	2.00	0.44
1:A:181:SER:O	1:A:183:PRO:HD3	2.18	0.44
1:A:114:GLU:HB3	1:A:132:ALA:HA	1.99	0.43
1:A:179:ILE:HA	1:A:180:PRO:HD3	1.91	0.43
1:B:240:VAL:HA	1:B:245:ASN:O	2.18	0.43
1:B:147:LEU:HD21	1:B:151:GLN:OE1	2.18	0.43
1:A:50:ALA:HB2	1:A:377:MET:HE2	2.01	0.43
1:A:50:ALA:HB2	1:A:377:MET:CE	2.49	0.43
1:B:243:GLU:O	1:B:243:GLU:HG2	2.19	0.43
1:A:45:LEU:HD22	1:A:368:PHE:HB2	2.01	0.43
1:A:123:LYS:HE3	1:A:245:ASN:CG	2.39	0.43
1:B:175:VAL:C	1:B:177:HIS:H	2.21	0.43
1:B:367:GLU:HG3	3:B:433:HOH:O	2.19	0.42
1:B:262:ARG:HH21	1:B:295:GLU:CD	2.21	0.42
1:B:66:LYS:HG3	1:B:391:THR:HG21	2.00	0.42
1:A:100:ASN:ND2	1:A:106:ASN:N	2.58	0.42
1:A:175:VAL:O	1:A:177:HIS:N	2.53	0.42
1:B:382:MET:HE3	1:B:389:ALA:HB2	1.99	0.42
1:A:265:GLN:HG3	1:A:297:LEU:HD13	2.01	0.42
1:B:396:LEU:CD1	1:B:402:ALA:HA	2.49	0.42
1:A:383:SER:HB3	1:A:386:ASP:CB	2.34	0.42
1:A:278:THR:CG2	1:A:282:ASN:O	2.68	0.42
1:A:43:PHE:HB3	1:A:240:VAL:HG12	2.01	0.42
1:B:82:ALA:O	1:B:86:VAL:HG22	2.20	0.41
1:A:145:ARG:HA	1:A:145:ARG:HD2	1.89	0.41
1:A:261:GLY:O	1:B:165:LEU:HA	2.20	0.41
1:B:323:MET:O	1:B:327:PHE:HB3	2.21	0.41
1:B:235:VAL:CG2	1:B:251:SER:HB3	2.50	0.41
1:B:284:TYR:HB3	1:B:347:ARG:HG3	2.01	0.41
1:B:156:SER:HB2	1:B:159:ARG:N	2.36	0.41
1:B:43:PHE:CG	1:B:44:VAL:N	2.88	0.41
1:A:76:LEU:HD23	1:A:76:LEU:HA	1.93	0.41
1:B:154:LYS:C	1:B:156:SER:H	2.24	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:124:SER:O	1:A:125:LEU:HB2	2.20	0.40
1:B:116:ASP:HB2	1:B:266:ALA:HB3	2.03	0.40
1:B:278:THR:HG22	1:B:284:TYR:C	2.41	0.40
1:A:239:VAL:O	1:A:246:VAL:HA	2.22	0.40
1:A:240:VAL:O	1:A:240:VAL:HG13	2.21	0.40
1:B:240:VAL:O	1:B:240:VAL:HG13	2.21	0.40
1:B:397:PRO:O	1:B:400:ALA:HB3	2.22	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	302/420 (72%)	275 (91%)	22 (7%)	5 (2%)	9 18
1	B	305/420 (73%)	271 (89%)	29 (10%)	5 (2%)	9 19
All	All	607/840 (72%)	546 (90%)	51 (8%)	10 (2%)	9 19

All (10) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	384	ALA
1	A	80	ALA
1	A	388	LYS
1	A	176	ASP
1	B	152	LYS
1	A	56	SER
1	B	266	ALA
1	B	293	CYS
1	A	154	LYS
1	B	387	GLY

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	233/331 (70%)	224 (96%)	9 (4%)	32	58
1	B	235/331 (71%)	225 (96%)	10 (4%)	29	54
All	All	468/662 (71%)	449 (96%)	19 (4%)	30	56

All (19) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	59	LYS
1	A	141	SER
1	A	172	ARG
1	A	265	GLN
1	A	277	ASN
1	A	302	LEU
1	A	304	ARG
1	A	313	GLU
1	A	395	ARG
1	B	71	LYS
1	B	86	VAL
1	B	126	ASN
1	B	145	ARG
1	B	168	GLU
1	B	265	GLN
1	B	313	GLU
1	B	329	SER
1	B	337	ASP
1	B	346	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (12) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	100	ASN
1	A	126	ASN
1	A	151	GLN

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Mol	Chain	Res	Type
1	A	277	ASN
1	B	70	GLN
1	B	100	ASN
1	B	106	ASN
1	B	138	ASN
1	B	144	ASN
1	B	265	GLN
1	B	277	ASN
1	B	317	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers (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)

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains (i)

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates (i)

EDS was not executed - this section is therefore empty.

6.4 Ligands (i)

EDS was not executed - this section is therefore empty.

6.5 Other polymers (i)

EDS was not executed - this section is therefore empty.