



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 01:21 PM GMT

PDB ID : 3TPT
Title : Structure of HipA(D309Q) bound to ADP
Authors : schumacher, M.A.; link, T.; Brennan, R.G.
Deposited on : 2011-09-08
Resolution : 2.25 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

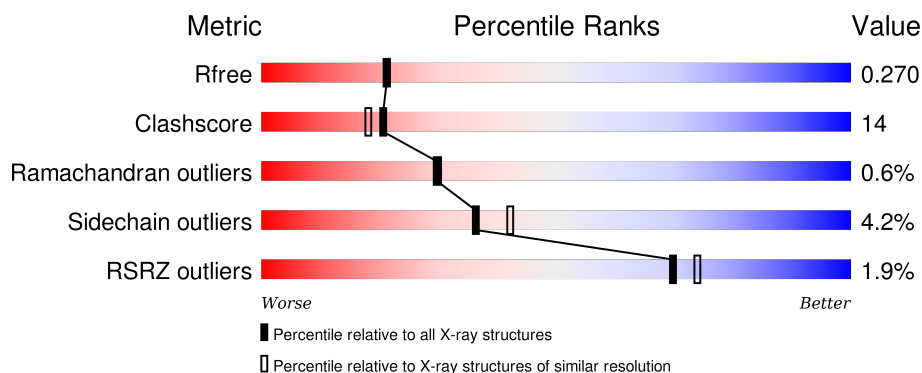
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 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	1640 (2.28-2.24)
Clashscore	102246	1095 (2.26-2.26)
Ramachandran outliers	100387	1063 (2.26-2.26)
Sidechain outliers	100360	1063 (2.26-2.26)
RSRZ outliers	91569	1647 (2.28-2.24)

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	440	
1	B	440	

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
3	MG	A	441	-	-	-	X

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6885 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.

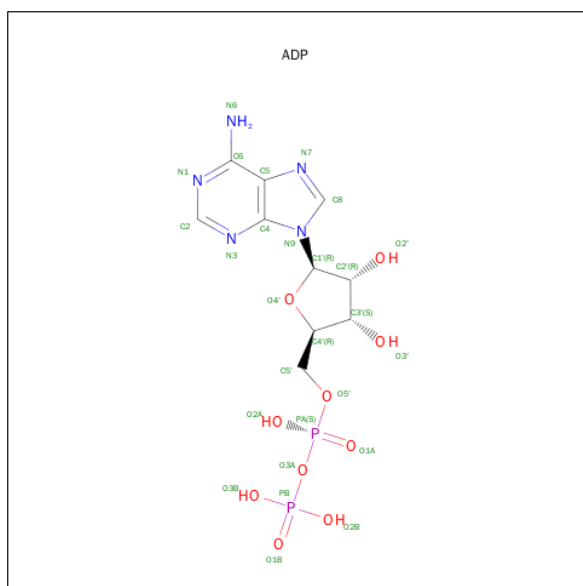
- Molecule 1 is a protein called Serine/threonine-protein kinase HipA.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	415	Total	C	N	O	S	0	0	0
			3286	2108	576	589	13			
1	B	419	Total	C	N	O	S	0	0	0
			3314	2121	582	598	13			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	243	ARG	GLU	CONFLICT	UNP P23874
A	309	GLN	ASP	ENGINEERED MUTATION	UNP P23874
B	243	ARG	GLU	CONFLICT	UNP P23874
B	309	GLN	ASP	ENGINEERED MUTATION	UNP P23874

- Molecule 2 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: $C_{10}H_{15}N_5O_{10}P_2$).

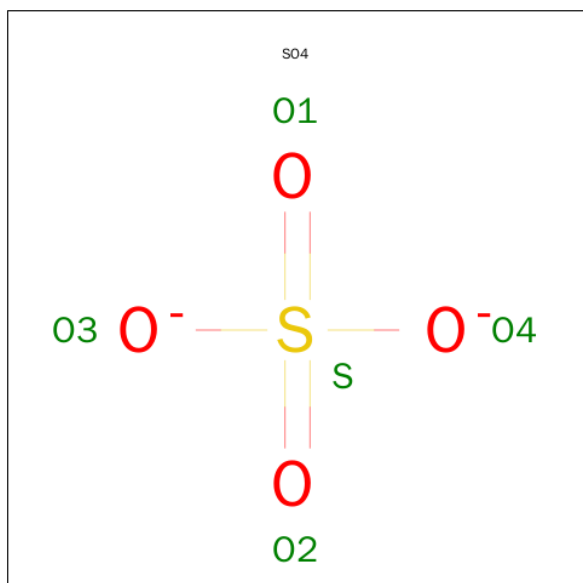


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			27	10	5	10	2		
2	B	1	Total	C	N	O	P	0	0
			27	10	5	10	2		

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	102	Total	O	0	0
			102	102		

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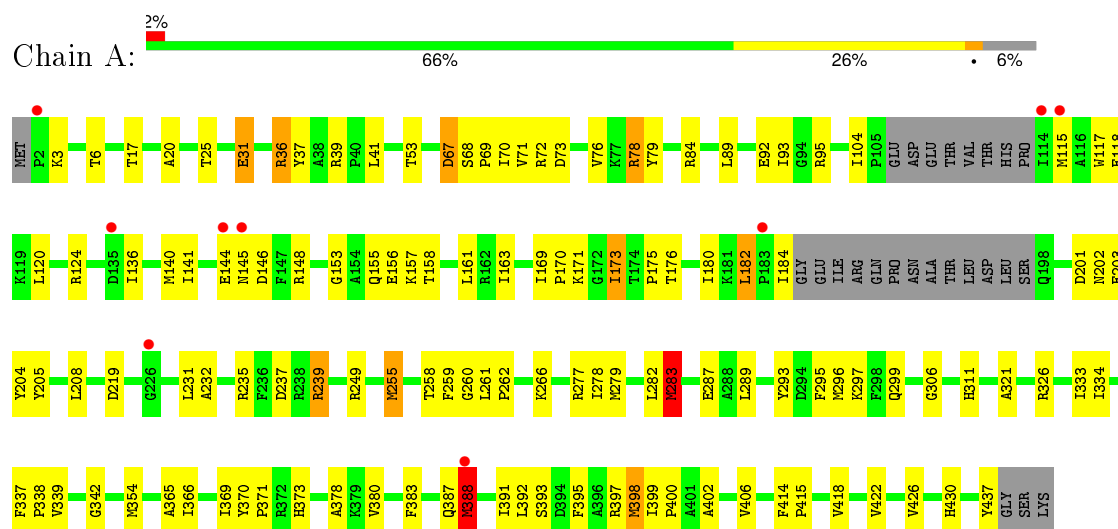
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	B	116	Total 116	O 116	0	0

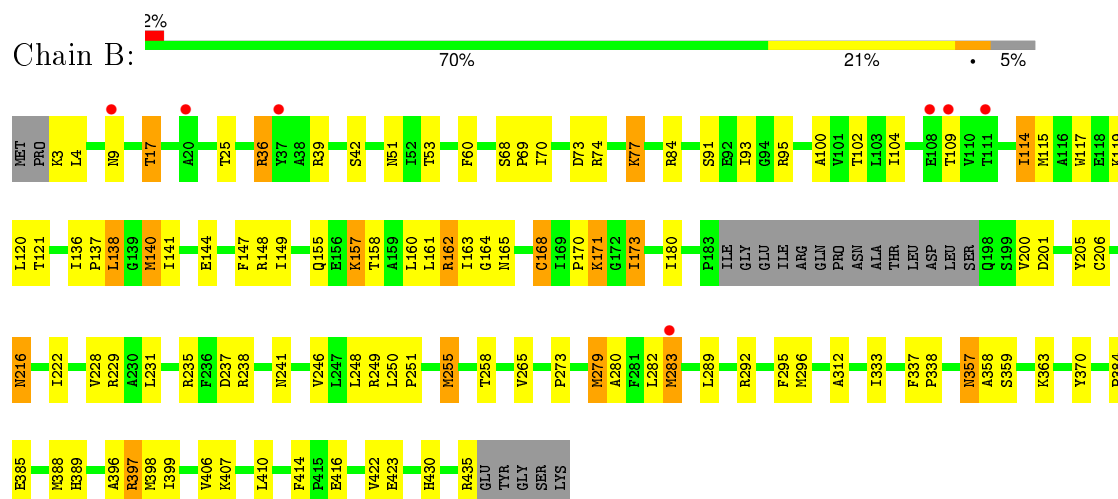
3 Residue-property plots

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: Serine/threonine-protein kinase HipA



• Molecule 1: Serine/threonine-protein kinase HipA



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	68.60Å 84.43Å 69.63Å 90.00° 91.66° 90.00°	Depositor
Resolution (Å)	25.20 – 2.25 25.20 – 2.25	Depositor EDS
% Data completeness (in resolution range)	96.9 (25.20-2.25) 93.3 (25.20-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.48 (at 2.26Å)	Xtriage
Refinement program	CNS 1.2	Depositor
R, R_{free}	0.232 , 0.267 0.240 , 0.270	Depositor DCC
R_{free} test set	1815 reflections (4.95%)	DCC
Wilson B-factor (Å ²)	42.0	Xtriage
Anisotropy	0.399	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 28.4	EDS
Estimated twinning fraction	0.028 for l,k,-h 0.043 for h,-k,-l 0.159 for l,-k,h	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.44$, $\langle L^2 \rangle = 0.27$	Xtriage
Outliers	0 of 36653 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6885	wwPDB-VP
Average B, all atoms (Å ²)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 38.00 % of the origin peak, indicating pseudo translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo translational symmetry is equal to 3.9386e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

¹ Intensities estimated from amplitudes.

² Theoretical 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: 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	0.49	9/3358 (0.3%)	0.60	0/4546
1	B	0.48	7/3387 (0.2%)	0.60	0/4589
All	All	0.48	16/6745 (0.2%)	0.60	0/9135

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	398	MET	CG-SD	6.48	1.98	1.81
1	A	279	MET	CG-SD	6.37	1.97	1.81
1	A	296	MET	CG-SD	5.86	1.96	1.81
1	B	255	MET	CG-SD	5.76	1.96	1.81
1	A	115	MET	CG-SD	5.74	1.96	1.81
1	A	255	MET	CG-SD	5.74	1.96	1.81
1	A	283	MET	CG-SD	5.72	1.96	1.81
1	A	398	MET	CG-SD	5.55	1.95	1.81
1	B	388	MET	CG-SD	5.50	1.95	1.81
1	B	279	MET	CG-SD	5.48	1.95	1.81
1	B	115	MET	CG-SD	5.39	1.95	1.81
1	A	354	MET	CG-SD	5.32	1.95	1.81
1	A	388	MET	CG-SD	5.26	1.94	1.81
1	B	296	MET	CG-SD	5.18	1.94	1.81
1	B	140	MET	CG-SD	5.14	1.94	1.81
1	A	140	MET	CG-SD	5.10	1.94	1.81

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts ⓘ

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3286	0	3342	108	0
1	B	3314	0	3362	87	0
2	A	27	0	12	0	0
2	B	27	0	12	1	0
3	A	1	0	0	0	0
3	B	2	0	0	0	0
4	A	5	0	0	0	0
4	B	5	0	0	0	0
5	A	102	0	0	4	0
5	B	116	0	0	2	0
All	All	6885	0	6728	194	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 14.

All (194) 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:279:MET:CE	1:B:282:LEU:HD12	1.91	0.99
1:B:279:MET:HE2	1:B:282:LEU:HD12	1.50	0.92
1:B:283:MET:HA	1:B:283:MET:HE3	1.52	0.92
1:A:173:ILE:HD13	1:A:173:ILE:H	1.38	0.87
1:B:358:ALA:HB2	1:B:363:LYS:HG3	1.56	0.86
1:A:136:ILE:HD13	1:A:141:ILE:HA	1.56	0.86
1:B:114:ILE:H	1:B:114:ILE:HD13	1.42	0.84
1:A:283:MET:HE3	1:A:283:MET:HA	1.65	0.79
1:B:407:LYS:HE2	1:B:423:GLU:OE1	1.83	0.79
1:A:202:ASN:ND2	1:A:338:PRO:HG3	1.98	0.78
1:A:370:TYR:H	1:A:373:HIS:HD2	1.32	0.78
1:B:161:LEU:HG	1:B:163:ILE:CD1	2.14	0.77
1:A:120:LEU:HD21	1:A:169:ILE:HG13	1.66	0.75
1:A:370:TYR:H	1:A:373:HIS:CD2	2.06	0.74
1:B:157:LYS:HB2	1:B:180:ILE:O	1.89	0.73
1:B:279:MET:HE1	1:B:295:PHE:HB3	1.70	0.73
1:B:283:MET:CE	1:B:283:MET:HA	2.17	0.73

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:25:THR:HG22	1:B:53:THR:HG22	1.72	0.71
1:A:25:THR:HG22	1:A:53:THR:HG22	1.73	0.71
1:B:173:ILE:H	1:B:173:ILE:HD13	1.56	0.70
1:A:36:ARG:HD3	1:A:36:ARG:H	1.55	0.70
1:A:93:ILE:HA	1:A:148:ARG:HB3	1.73	0.69
1:A:283:MET:CE	1:A:283:MET:HA	2.22	0.69
1:A:255:MET:O	1:A:258:THR:HG22	1.93	0.68
1:B:36:ARG:HD3	1:B:36:ARG:N	2.09	0.68
1:B:222:ILE:HD13	1:B:229:ARG:NH1	2.08	0.68
1:A:239:ARG:HH11	1:A:239:ARG:HB3	1.59	0.68
1:A:205:TYR:CE2	1:A:422:VAL:HG11	2.28	0.67
1:B:36:ARG:HD3	1:B:36:ARG:H	1.57	0.67
1:B:161:LEU:HG	1:B:163:ILE:HD11	1.75	0.67
1:A:3:LYS:HG3	1:A:17:THR:HG22	1.79	0.65
1:B:255:MET:O	1:B:258:THR:HG22	1.97	0.65
1:A:36:ARG:HB2	1:A:37:TYR:CD1	2.33	0.64
1:B:157:LYS:NZ	2:B:501:ADP:O1B	2.27	0.63
1:B:3:LYS:HD3	1:B:17:THR:HB	1.79	0.63
1:B:357:ASN:N	1:B:357:ASN:HD22	1.97	0.63
1:A:17:THR:HB	5:A:481:HOH:O	1.99	0.62
1:A:289:LEU:H	1:A:289:LEU:HD22	1.65	0.61
1:B:17:THR:HG23	1:B:25:THR:OG1	2.00	0.61
1:B:222:ILE:HD13	1:B:229:ARG:HH11	1.67	0.59
1:B:36:ARG:CD	1:B:36:ARG:H	2.15	0.59
1:A:378:ALA:HB2	1:A:388:MET:HG3	1.84	0.58
1:A:79:TYR:HB3	1:A:92:GLU:OE1	2.04	0.57
1:B:25:THR:CG2	1:B:53:THR:HG22	2.35	0.57
1:A:36:ARG:H	1:A:36:ARG:CD	2.16	0.56
1:B:36:ARG:HH11	1:B:36:ARG:H	1.52	0.56
1:A:392:LEU:HD13	1:A:437:TYR:OH	2.06	0.56
1:B:114:ILE:H	1:B:114:ILE:CD1	2.16	0.56
1:A:37:TYR:HB3	1:A:321:ALA:HB1	1.87	0.56
1:B:385:GLU:HG2	1:B:389:HIS:CD2	2.41	0.56
1:B:136:ILE:HG23	1:B:137:PRO:HD2	1.88	0.55
1:A:422:VAL:O	1:A:426:VAL:HG23	2.07	0.55
1:A:31:GLU:OE1	1:A:31:GLU:N	2.40	0.55
1:A:180:ILE:HD13	1:A:232:ALA:HA	1.90	0.54
1:B:358:ALA:CB	1:B:363:LYS:HG3	2.34	0.54
1:B:201:ASP:HB3	1:B:414:PHE:CE2	2.43	0.54
1:B:160:LEU:HD12	1:B:180:ILE:CD1	2.38	0.54
1:B:283:MET:HE3	1:B:292:ARG:HD2	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:205:TYR:HE2	1:A:422:VAL:HG11	1.73	0.54
1:A:78:ARG:O	1:A:78:ARG:HD2	2.07	0.54
1:B:396:ALA:HB1	1:B:435:ARG:HB3	1.90	0.54
1:B:138:LEU:HD12	1:B:140:MET:HE1	1.89	0.54
1:A:158:THR:HG22	1:A:180:ILE:HB	1.90	0.53
1:A:378:ALA:CB	1:A:388:MET:HG3	2.37	0.53
1:A:161:LEU:HD21	1:A:163:ILE:HD11	1.90	0.53
1:B:170:PRO:C	1:B:171:LYS:HD2	2.28	0.53
1:A:118:GLU:CD	1:A:124:ARG:HH12	2.10	0.53
1:A:337:PHE:N	1:A:338:PRO:HD2	2.24	0.53
1:B:255:MET:HA	1:B:258:THR:HG22	1.89	0.53
1:A:95:ARG:NE	1:A:104:ILE:HG22	2.24	0.53
1:A:295:PHE:O	1:A:299:GLN:HG2	2.09	0.52
1:B:399:ILE:HB	1:B:430:HIS:CD2	2.44	0.52
1:A:20:ALA:O	1:B:289:LEU:HD12	2.10	0.52
1:A:173:ILE:CD1	1:A:173:ILE:H	2.17	0.52
1:A:182:LEU:HD22	1:A:182:LEU:N	2.25	0.51
1:B:70:ILE:O	1:B:73:ASP:HB2	2.10	0.51
1:B:157:LYS:HG2	1:B:158:THR:N	2.24	0.51
1:A:365:ALA:O	1:A:369:ILE:HG13	2.11	0.51
1:A:120:LEU:HD21	1:A:169:ILE:CG1	2.38	0.51
1:A:399:ILE:HB	1:A:430:HIS:CD2	2.45	0.51
1:B:161:LEU:HG	1:B:163:ILE:HD12	1.92	0.50
1:A:204:TYR:CE2	1:A:208:LEU:HD11	2.47	0.50
1:B:163:ILE:HD13	1:B:168:CYS:SG	2.52	0.50
1:B:119:LYS:HG2	1:B:120:LEU:N	2.26	0.50
1:B:141:ILE:HD12	1:B:141:ILE:O	2.12	0.49
1:B:117:TRP:CE2	1:B:163:ILE:HG12	2.47	0.49
1:B:163:ILE:N	1:B:163:ILE:HD12	2.28	0.49
1:A:158:THR:CG2	1:A:180:ILE:HB	2.42	0.49
1:A:182:LEU:HD22	1:A:182:LEU:H	1.77	0.49
1:A:260:GLY:HA2	5:A:523:HOH:O	2.12	0.49
1:A:117:TRP:HA	1:A:171:LYS:HG2	1.94	0.49
1:A:297:LYS:HB2	1:A:391:ILE:HD12	1.95	0.49
1:A:72:ARG:O	1:A:76:VAL:HG23	2.12	0.49
1:B:42:SER:HB2	1:B:60:PHE:CE1	2.47	0.49
1:A:289:LEU:N	1:A:289:LEU:HD22	2.28	0.49
1:A:387:GLN:O	1:A:391:ILE:HG12	2.13	0.48
1:A:333:ILE:O	1:A:334:ILE:HD13	2.13	0.48
1:A:180:ILE:HA	1:A:231:LEU:O	2.14	0.48
1:A:380:VAL:HG12	1:A:380:VAL:O	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:161:LEU:CG	1:B:163:ILE:HD11	2.42	0.48
1:B:337:PHE:HB2	1:B:338:PRO:HD3	1.96	0.48
1:A:170:PRO:HG3	1:A:176:THR:HG22	1.96	0.48
1:A:235:ARG:HB3	1:A:237:ASP:OD1	2.13	0.48
1:A:418:VAL:O	1:A:422:VAL:HG23	2.14	0.47
1:A:76:VAL:HG22	1:A:89:LEU:HD21	1.95	0.47
1:B:206:CYS:HB3	1:B:333:ILE:HG22	1.97	0.47
1:B:238:ARG:HA	1:B:248:LEU:O	2.14	0.47
1:B:77:LYS:HD2	1:B:77:LYS:C	2.35	0.47
1:A:3:LYS:HG3	1:A:17:THR:CG2	2.44	0.46
1:A:297:LYS:HG3	1:A:391:ILE:HG23	1.98	0.46
1:A:266:LYS:HD2	1:A:311:HIS:CE1	2.50	0.46
1:B:273:PRO:HG2	1:B:312:ALA:HB2	1.97	0.46
1:B:170:PRO:O	1:B:171:LYS:HD2	2.16	0.46
1:B:205:TYR:HE2	1:B:422:VAL:HG11	1.81	0.46
1:A:393:SER:HB3	1:A:437:TYR:CD2	2.51	0.46
1:A:68:SER:O	1:A:72:ARG:HG3	2.15	0.45
1:B:406:VAL:O	1:B:410:LEU:HG	2.16	0.45
1:A:395:PHE:O	1:A:399:ILE:HG12	2.17	0.45
1:A:339:VAL:O	1:A:339:VAL:HG12	2.16	0.45
1:A:414:PHE:CG	1:A:415:PRO:HD2	2.52	0.45
1:A:202:ASN:ND2	1:A:338:PRO:CG	2.77	0.45
1:A:278:ILE:O	1:A:282:LEU:HG	2.17	0.45
1:A:293:TYR:CZ	1:A:391:ILE:HD11	2.53	0.44
1:A:70:ILE:O	1:A:73:ASP:HB2	2.17	0.44
1:A:156:GLU:H	1:A:184:ILE:HD12	1.83	0.44
1:A:175:PRO:HG3	1:A:249:ARG:CZ	2.48	0.44
1:B:358:ALA:O	1:B:359:SER:C	2.56	0.44
1:A:25:THR:CG2	1:A:53:THR:HG22	2.44	0.44
1:A:155:GLN:CG	1:A:184:ILE:HD11	2.48	0.44
1:B:241:ASN:ND2	1:B:246:VAL:HB	2.33	0.44
1:A:117:TRP:CE2	1:A:163:ILE:HD12	2.52	0.44
1:A:399:ILE:HB	1:A:400:PRO:HD3	2.00	0.44
1:B:162:ARG:HD3	1:B:164:GLY:O	2.17	0.44
1:A:326:ARG:HG2	5:A:451:HOH:O	2.17	0.43
1:A:202:ASN:HD21	1:A:338:PRO:HG3	1.76	0.43
1:A:383:PHE:CE2	1:A:388:MET:HG2	2.52	0.43
1:A:157:LYS:HA	1:A:182:LEU:HD22	2.00	0.43
1:A:397:ARG:HH11	1:A:397:ARG:HG2	1.83	0.43
1:A:67:ASP:OD1	1:A:153:GLY:N	2.49	0.43
1:B:397:ARG:HG2	1:B:397:ARG:HH11	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:383:PHE:CD2	1:A:388:MET:HG2	2.53	0.43
1:A:144:GLU:HG3	1:A:146:ASP:H	1.84	0.43
1:B:235:ARG:HB3	1:B:237:ASP:OD1	2.18	0.43
1:A:6:THR:HG21	1:A:41:LEU:HD11	2.00	0.43
1:A:120:LEU:HD21	1:A:169:ILE:CD1	2.48	0.43
1:A:201:ASP:HB3	1:A:414:PHE:CZ	2.54	0.43
1:A:68:SER:HB3	1:A:71:VAL:HG23	1.99	0.43
1:B:74:ARG:O	1:B:77:LYS:HG3	2.19	0.43
1:B:155:GLN:HE21	1:B:155:GLN:HB3	1.59	0.43
1:A:78:ARG:HD2	1:A:78:ARG:C	2.39	0.43
1:A:41:LEU:HD23	1:A:41:LEU:HA	1.85	0.43
1:B:95:ARG:NH2	1:B:148:ARG:HG3	2.33	0.43
1:B:370:TYR:HD2	5:B:474:HOH:O	2.00	0.43
1:B:205:TYR:CE2	1:B:422:VAL:HG11	2.53	0.43
1:B:397:ARG:HG2	1:B:397:ARG:NH1	2.33	0.43
1:A:144:GLU:HG3	1:A:145:ASN:N	2.34	0.42
1:B:280:ALA:O	1:B:283:MET:HB2	2.19	0.42
1:B:200:VAL:HG13	1:B:231:LEU:HB2	2.01	0.42
1:B:279:MET:HE1	1:B:295:PHE:CB	2.45	0.42
1:A:76:VAL:HG21	1:A:84:ARG:HG2	2.01	0.42
1:A:306:GLY:HA2	1:A:366:ILE:HD13	2.00	0.42
1:A:261:LEU:HA	1:A:262:PRO:HD3	1.85	0.42
1:A:255:MET:HA	1:A:258:THR:HG22	2.01	0.42
1:A:201:ASP:HB3	1:A:414:PHE:CE2	2.55	0.42
1:A:370:TYR:HB3	1:A:371:PRO:HD2	2.01	0.42
1:A:402:ALA:O	1:A:406:VAL:HG23	2.20	0.42
1:A:78:ARG:HB3	1:A:79:TYR:CD1	2.55	0.41
1:A:287:GLU:OE2	1:A:287:GLU:HA	2.20	0.41
1:A:369:ILE:HA	1:A:373:HIS:CD2	2.55	0.41
1:B:36:ARG:CD	1:B:36:ARG:N	2.78	0.41
1:B:95:ARG:O	1:B:102:THR:HG23	2.20	0.41
1:B:84:ARG:NH2	5:B:506:HOH:O	2.54	0.41
1:B:141:ILE:HD11	1:B:144:GLU:CG	2.50	0.41
1:B:4:LEU:HD21	1:B:91:SER:HA	2.02	0.41
1:B:68:SER:HA	1:B:69:PRO:HD3	1.94	0.41
1:B:95:ARG:HG3	1:B:104:ILE:CD1	2.51	0.41
1:B:100:ALA:HA	1:B:249:ARG:HB3	2.02	0.41
1:B:157:LYS:CG	1:B:158:THR:N	2.83	0.41
1:A:255:MET:HG3	1:A:278:ILE:HD12	2.03	0.41
1:B:228:VAL:HG22	1:B:229:ARG:N	2.36	0.41
1:A:203:GLU:HB2	1:A:334:ILE:CD1	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:155:GLN:HG3	1:A:184:ILE:HD11	2.03	0.41
1:A:397:ARG:HG3	1:A:398:MET:N	2.35	0.41
1:B:93:ILE:HB	1:B:149:ILE:O	2.20	0.41
1:A:68:SER:HA	1:A:69:PRO:HD3	1.94	0.41
1:B:93:ILE:HA	1:B:148:ARG:HB3	2.02	0.41
1:A:255:MET:O	1:A:259:PHE:HD1	2.04	0.40
1:B:216:ASN:HD22	1:B:216:ASN:HA	1.67	0.40
1:B:171:LYS:N	1:B:171:LYS:HD2	2.36	0.40
1:B:250:LEU:HA	1:B:251:PRO:HD3	1.90	0.40
1:B:265:VAL:O	1:B:265:VAL:HG23	2.21	0.40
1:A:437:TYR:CD1	1:A:437:TYR:N	2.89	0.40
1:A:277:ARG:HG3	5:A:446:HOH:O	2.21	0.40
1:B:161:LEU:CD2	1:B:163:ILE:HD11	2.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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	409/440 (93%)	386 (94%)	22 (5%)	1 (0%)	52	61
1	B	415/440 (94%)	392 (94%)	19 (5%)	4 (1%)	19	16
All	All	824/880 (94%)	778 (94%)	41 (5%)	5 (1%)	30	30

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	109	THR
1	B	9	ASN
1	B	138	LEU
1	A	342	GLY
1	B	384	PRO

5.3.2 Protein sidechains ⓘ

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	352/374 (94%)	341 (97%)	11 (3%)	47	58
1	B	356/374 (95%)	337 (95%)	19 (5%)	28	30
All	All	708/748 (95%)	678 (96%)	30 (4%)	36	42

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	31	GLU
1	A	36	ARG
1	A	39	ARG
1	A	67	ASP
1	A	78	ARG
1	A	173	ILE
1	A	182	LEU
1	A	219	ASP
1	A	239	ARG
1	A	283	MET
1	A	388	MET
1	B	17	THR
1	B	36	ARG
1	B	39	ARG
1	B	51	ASN
1	B	77	LYS
1	B	114	ILE
1	B	121	THR
1	B	147	PHE
1	B	157	LYS
1	B	162	ARG
1	B	165	ASN
1	B	168	CYS
1	B	171	LYS
1	B	173	ILE
1	B	216	ASN
1	B	283	MET

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Mol	Chain	Res	Type
1	B	357	ASN
1	B	397	ARG
1	B	416	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (11) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	155	GLN
1	A	216	ASN
1	A	309	GLN
1	A	320	GLN
1	A	357	ASN
1	A	373	HIS
1	B	51	ASN
1	B	155	GLN
1	B	216	ASN
1	B	320	GLN
1	B	357	ASN

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 7 ligands modelled in this entry, 3 are monoatomic - leaving 4 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length

(or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	ADP	A	500	3	22,29,29	1.97	6 (27%)	27,45,45	2.36	9 (33%)
4	SO4	A	833	-	4,4,4	0.21	0	6,6,6	0.12	0
2	ADP	B	501	3	22,29,29	1.95	6 (27%)	27,45,45	2.33	9 (33%)
4	SO4	B	834	-	4,4,4	0.22	0	6,6,6	0.10	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	ADP	A	500	3	-	0/12/32/32	0/3/3/3
4	SO4	A	833	-	-	0/0/0/0	0/0/0/0
2	ADP	B	501	3	-	0/12/32/32	0/3/3/3
4	SO4	B	834	-	-	0/0/0/0	0/0/0/0

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	ADP	PB-O2B	-2.43	1.46	1.54
2	A	500	ADP	PB-O2B	-2.32	1.46	1.54
2	A	500	ADP	C2-N1	2.00	1.37	1.33
2	B	501	ADP	C2-N1	2.03	1.37	1.33
2	A	500	ADP	O4'-C1'	2.78	1.44	1.41
2	B	501	ADP	PB-O3B	2.83	1.64	1.54
2	A	500	ADP	PB-O3B	3.11	1.65	1.54
2	B	501	ADP	O4'-C1'	3.44	1.45	1.41
2	A	500	ADP	C2-N3	3.74	1.38	1.32
2	B	501	ADP	C2-N3	4.12	1.39	1.32
2	B	501	ADP	C4-N3	4.75	1.42	1.35
2	A	500	ADP	C4-N3	5.35	1.43	1.35

All (18) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	500	ADP	N3-C2-N1	-4.48	125.47	128.89
2	B	501	ADP	N3-C2-N1	-4.45	125.49	128.89
2	A	500	ADP	C1'-N9-C4	-3.48	121.70	126.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	ADP	O2B-PB-O3A	-2.78	92.49	105.09
2	B	501	ADP	C1'-N9-C4	-2.69	122.88	126.94
2	A	500	ADP	O3B-PB-O3A	-2.59	93.32	105.09
2	A	500	ADP	O2B-PB-O3A	-2.43	94.08	105.09
2	B	501	ADP	O3B-PB-O3A	-2.30	94.64	105.09
2	B	501	ADP	PA-O3A-PB	2.39	140.67	132.67
2	B	501	ADP	C4-C5-N7	2.39	111.68	109.48
2	A	500	ADP	PA-O3A-PB	2.43	140.80	132.67
2	A	500	ADP	C4-C5-N7	2.60	111.87	109.48
2	A	500	ADP	C4'-O4'-C1'	2.97	112.98	109.72
2	B	501	ADP	C4'-O4'-C1'	3.22	113.26	109.72
2	B	501	ADP	C2'-C1'-N9	3.43	119.54	114.29
2	A	500	ADP	C2'-C1'-N9	3.77	120.05	114.29
2	A	500	ADP	O2B-PB-O1B	7.42	134.48	110.58
2	B	501	ADP	O2B-PB-O1B	7.60	135.06	110.58

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	501	ADP	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [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	415/440 (94%)	-0.06	9 (2%) 65 70	26, 42, 67, 85	0
1	B	419/440 (95%)	-0.15	7 (1%) 73 77	27, 40, 63, 84	0
All	All	834/880 (94%)	-0.10	16 (1%) 70 74	26, 41, 66, 85	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	114	ILE	7.7
1	B	20	ALA	4.0
1	A	145	ASN	4.0
1	B	109	THR	3.5
1	A	183	PRO	3.4
1	B	37	TYR	3.2
1	A	135	ASP	3.1
1	B	9	ASN	3.0
1	B	108	GLU	2.9
1	A	2	PRO	2.8
1	B	111	THR	2.7
1	A	115	MET	2.5
1	A	226	GLY	2.3
1	B	283	MET	2.2
1	A	388	MET	2.1
1	A	144	GLU	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(\AA^2)	Q<0.9
3	MG	A	441	1/1	0.71	0.15	2.45	60,60,60,60	0
4	SO4	B	834	5/5	0.92	0.14	-0.30	79,79,81,82	0
2	ADP	A	500	27/27	0.94	0.11	-0.32	39,47,52,55	0
4	SO4	A	833	5/5	0.97	0.12	-0.34	65,66,67,69	0
2	ADP	B	501	27/27	0.96	0.10	-0.50	27,36,48,51	0
3	MG	B	442	1/1	0.77	0.16	-	52,52,52,52	0
3	MG	B	441	1/1	0.97	0.18	-	59,59,59,59	0

6.5 Other polymers [i](#)

There are no such residues in this entry.