



Full wwPDB X-ray Structure Validation Report ⓘ

May 26, 2020 – 02:12 pm BST

PDB ID : 6ACX
Title : Crystal structure of Mycobacterium smegmatis Mfd in complex with ADP + Pi at 3.5 Å resolution.
Authors : Putta, S.; Fox, G.C.; Walsh, M.A.; Rao, D.N.; Nagaraja, V.; Natesh, R.
Deposited on : 2018-07-27
Resolution : 3.50 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

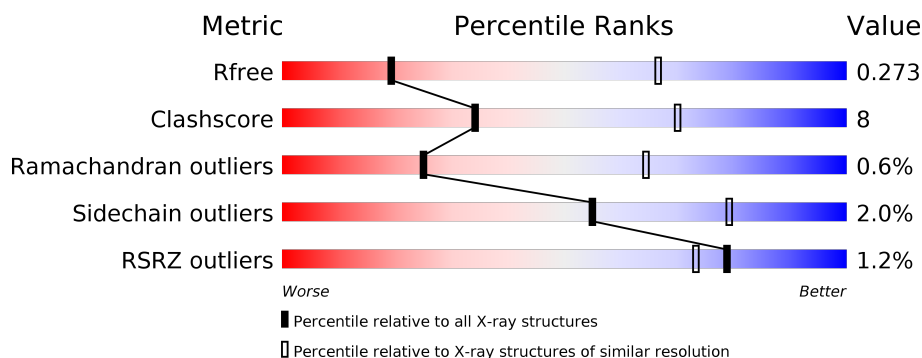
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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}	130704	1659 (3.60-3.40)
Clashscore	141614	1036 (3.58-3.42)
Ramachandran outliers	138981	1005 (3.58-3.42)
Sidechain outliers	138945	1006 (3.58-3.42)
RSRZ outliers	127900	1559 (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 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 $\leq 5\%$. 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	1235	<div> <div>%</div> <div> <div></div> <div>74%</div> <div>20%</div> <div>• 5%</div> </div> </div>
1	B	1235	<div> <div>%</div> <div> <div></div> <div>76%</div> <div>18%</div> <div>• 5%</div> </div> </div>

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 17349 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 Mycobacterium smegmatis Mfd.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	1175	Total	C	N	O	S	0	0	0
			8626	5416	1539	1641	30			
1	B	1168	Total	C	N	O	S	0	0	0
			8539	5362	1515	1631	31			

There are 40 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP I7G7M2
A	-18	GLY	-	expression tag	UNP I7G7M2
A	-17	SER	-	expression tag	UNP I7G7M2
A	-16	SER	-	expression tag	UNP I7G7M2
A	-15	HIS	-	expression tag	UNP I7G7M2
A	-14	HIS	-	expression tag	UNP I7G7M2
A	-13	HIS	-	expression tag	UNP I7G7M2
A	-12	HIS	-	expression tag	UNP I7G7M2
A	-11	HIS	-	expression tag	UNP I7G7M2
A	-10	HIS	-	expression tag	UNP I7G7M2
A	-9	SER	-	expression tag	UNP I7G7M2
A	-8	SER	-	expression tag	UNP I7G7M2
A	-7	GLY	-	expression tag	UNP I7G7M2
A	-6	LEU	-	expression tag	UNP I7G7M2
A	-5	VAL	-	expression tag	UNP I7G7M2
A	-4	PRO	-	expression tag	UNP I7G7M2
A	-3	ARG	-	expression tag	UNP I7G7M2
A	-2	GLY	-	expression tag	UNP I7G7M2
A	-1	SER	-	expression tag	UNP I7G7M2
A	0	HIS	-	expression tag	UNP I7G7M2
B	-19	MET	-	initiating methionine	UNP I7G7M2
B	-18	GLY	-	expression tag	UNP I7G7M2
B	-17	SER	-	expression tag	UNP I7G7M2
B	-16	SER	-	expression tag	UNP I7G7M2
B	-15	HIS	-	expression tag	UNP I7G7M2

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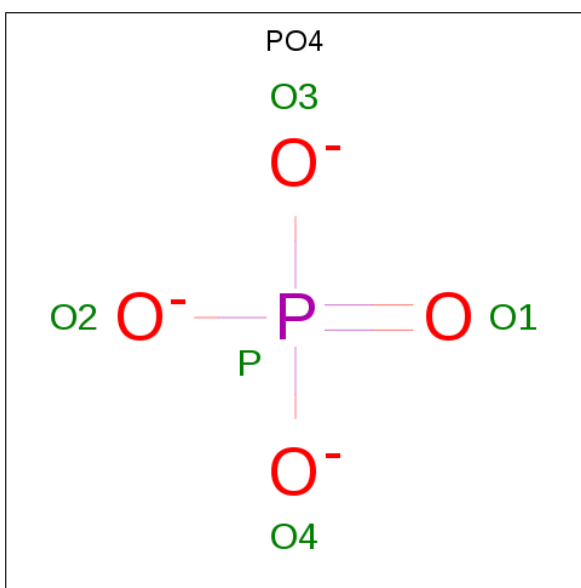
Chain	Residue	Modelled	Actual	Comment	Reference
B	-14	HIS	-	expression tag	UNP I7G7M2
B	-13	HIS	-	expression tag	UNP I7G7M2
B	-12	HIS	-	expression tag	UNP I7G7M2
B	-11	HIS	-	expression tag	UNP I7G7M2
B	-10	HIS	-	expression tag	UNP I7G7M2
B	-9	SER	-	expression tag	UNP I7G7M2
B	-8	SER	-	expression tag	UNP I7G7M2
B	-7	GLY	-	expression tag	UNP I7G7M2
B	-6	LEU	-	expression tag	UNP I7G7M2
B	-5	VAL	-	expression tag	UNP I7G7M2
B	-4	PRO	-	expression tag	UNP I7G7M2
B	-3	ARG	-	expression tag	UNP I7G7M2
B	-2	GLY	-	expression tag	UNP I7G7M2
B	-1	SER	-	expression tag	UNP I7G7M2
B	0	HIS	-	expression tag	UNP I7G7M2

- Molecule 2 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



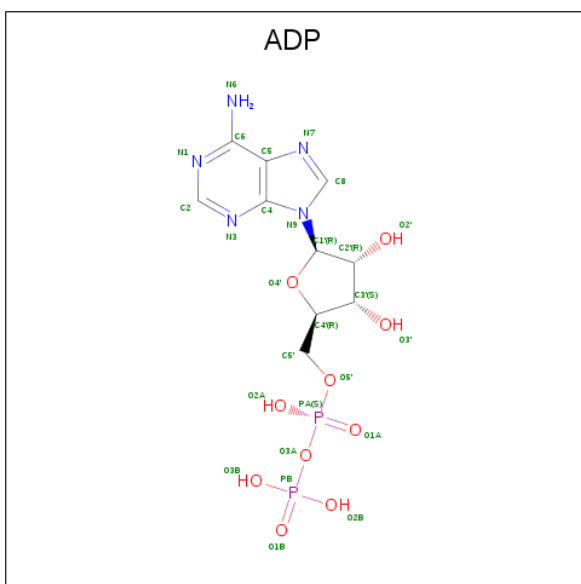
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	A	1	Total	O	S	0	0
			5	4	1		
2	B	1	Total	O	S	0	0
			5	4	1		

- Molecule 3 is PHOSPHATE ION (three-letter code: PO4) (formula: O₄P).



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

- Molecule 4 is ADENOSINE-5'-DIPHOSPHATE (three-letter code: ADP) (formula: C₁₀H₁₅N₅O₁₀P₂).



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

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

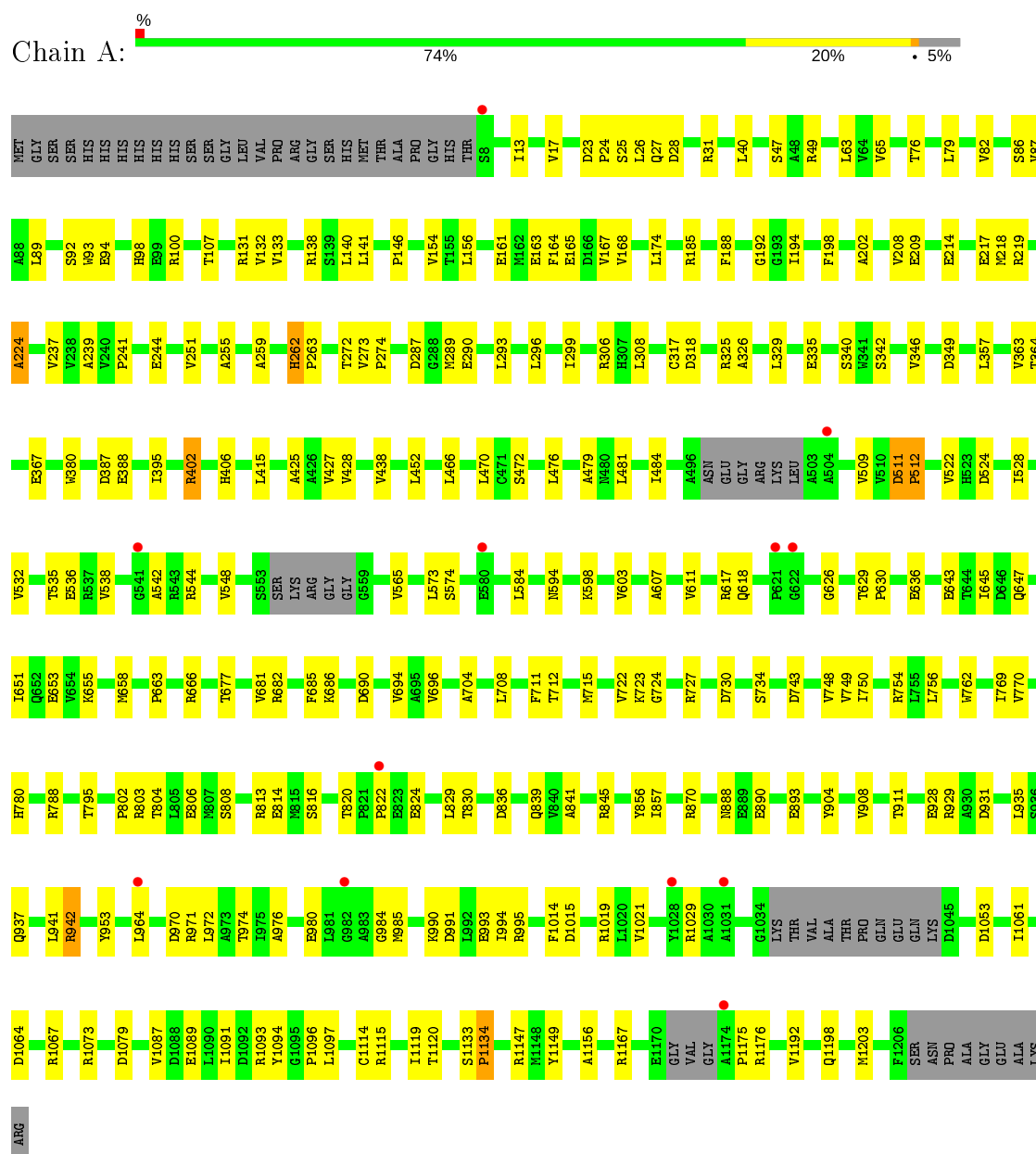
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	54	Total	O	0	0
			54	54		
5	B	51	Total	O	0	0
			51	51		

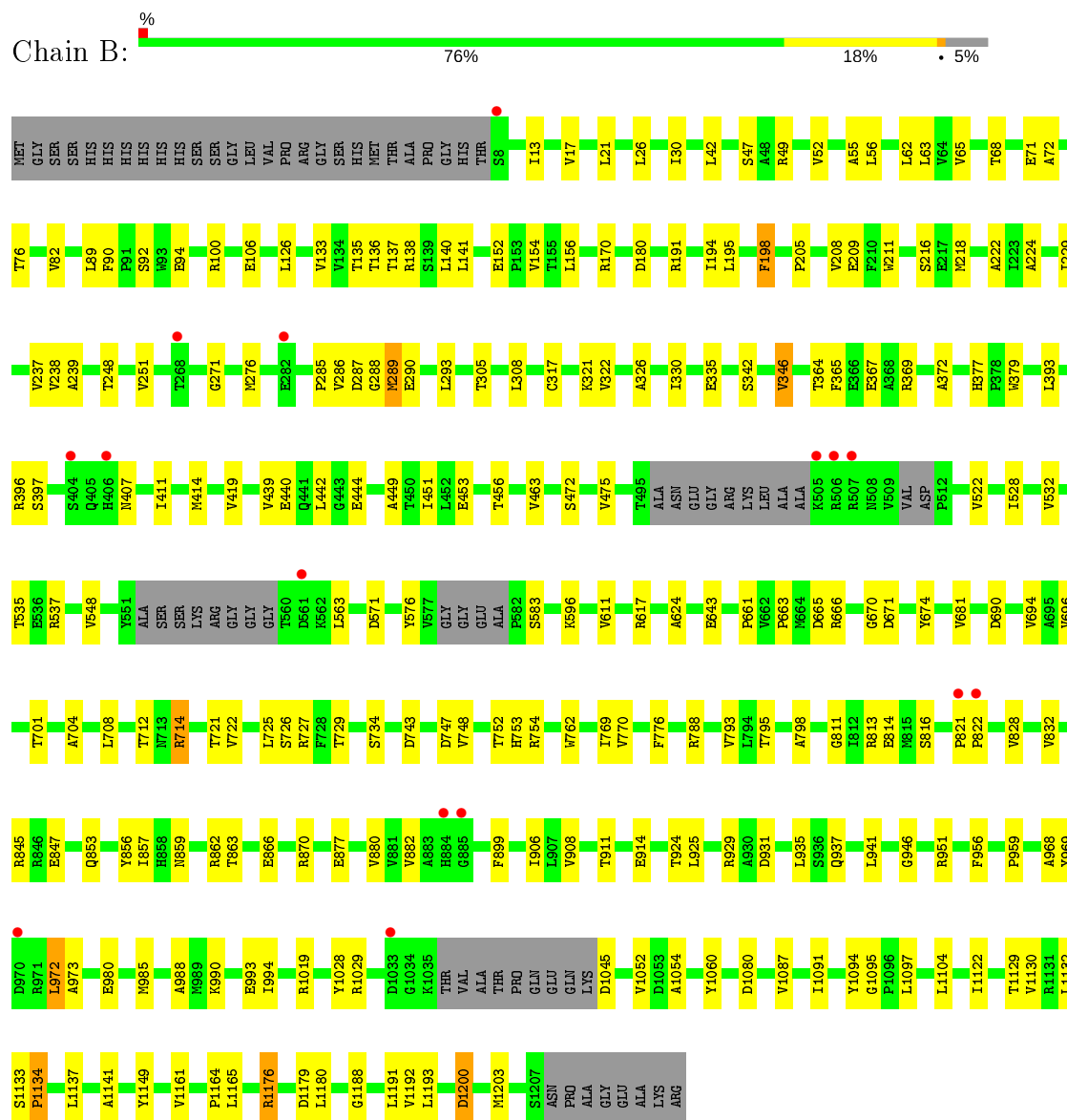
3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes 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: *Mycobacterium smegmatis* Mfd



- Molecule 1: *Mycobacterium smegmatis* Mfd



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	83.66Å 160.44Å 214.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.10 – 3.50 31.50 – 3.50	Depositor EDS
% Data completeness (in resolution range)	96.6 (30.10-3.50) 96.7 (31.50-3.50)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	0.22	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.95 (at 3.47Å)	Xtriage
Refinement program	PHENIX (1.11.1 _2575: 000)	Depositor
R, R_{free}	0.226 , 0.273 0.226 , 0.273	Depositor DCC
R_{free} test set	1770 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å ²)	66.1	Xtriage
Anisotropy	0.149	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.23 , 37.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.38$, $\langle L^2 \rangle = 0.21$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.86	EDS
Total number of atoms	17349	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.88% 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, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PO4, 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	0.25	0/8781	0.43	0/11992
1	B	0.24	0/8691	0.42	0/11874
All	All	0.24	0/17472	0.43	0/23866

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

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	8626	0	8374	148	0
1	B	8539	0	8256	140	0
2	A	15	0	0	3	0
2	B	5	0	0	0	0
3	A	5	0	0	0	0
4	A	27	0	12	3	0
4	B	27	0	12	2	0
5	A	54	0	0	1	0
5	B	51	0	0	1	0
All	All	17349	0	16654	286	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 8.

All (286) 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:511:ASP:H	1:A:512:PRO:HD3	1.36	0.89
1:A:617:ARG:HG3	1:A:663:PRO:HG2	1.67	0.76
1:B:968:ALA:O	1:B:972:LEU:HB2	1.87	0.75
1:B:832:VAL:HG12	1:B:956:PHE:HB2	1.72	0.72
1:A:251:VAL:HG23	1:A:296:LEU:HB3	1.71	0.71
1:B:674:TYR:HA	1:B:822:PRO:HG3	1.73	0.70
1:A:723:LYS:O	1:A:749:VAL:HA	1.93	0.69
1:B:929:ARG:HG2	1:B:959:PRO:HD3	1.73	0.69
1:A:273:VAL:HG23	1:A:274:PRO:HD3	1.75	0.69
1:A:804:THR:O	1:A:808:SER:HB3	1.93	0.68
1:A:857:ILE:HD12	1:A:941:LEU:HD13	1.76	0.68
1:B:704:ALA:O	1:B:708:LEU:HB2	1.94	0.67
1:A:788:ARG:HH21	1:A:813:ARG:HH12	1.43	0.67
1:A:658:MET:HE3	1:A:686:LYS:HD3	1.76	0.67
1:B:734:SER:OG	1:B:754:ARG:NH2	2.27	0.66
1:B:535:THR:OG1	1:B:537:ARG:NH2	2.28	0.66
1:A:255:ALA:O	1:A:259:ALA:HB2	1.96	0.66
1:B:548:VAL:HA	1:B:563:LEU:O	1.96	0.66
1:B:211:TRP:HB2	1:B:216:SER:HB3	1.77	0.65
1:B:694:VAL:HG22	1:B:748:VAL:HG12	1.78	0.65
1:A:904:TYR:HE1	1:B:877:GLU:HB3	1.62	0.65
1:A:208:VAL:HG12	1:A:218:MET:HG2	1.79	0.64
1:A:98:HIS:NE2	1:A:357:LEU:O	2.30	0.64
1:B:522:VAL:HG23	1:B:576:TYR:HB2	1.80	0.64
1:A:822:PRO:HB2	4:A:1305:ADP:C6	2.33	0.64
1:B:222:ALA:HB2	1:B:229:ILE:HD11	1.80	0.64
1:A:511:ASP:H	1:A:512:PRO:CD	2.10	0.63
1:A:532:VAL:HB	1:A:548:VAL:HG23	1.80	0.63
1:B:532:VAL:HB	1:B:548:VAL:HG23	1.79	0.63
1:A:194:ILE:HG12	1:A:209:GLU:HG2	1.79	0.63
1:A:156:LEU:HB2	1:A:237:VAL:HG22	1.81	0.63
1:A:629:THR:HB	1:A:630:PRO:HD2	1.80	0.63
1:B:617:ARG:NH1	1:B:663:PRO:O	2.32	0.62
1:B:980:GLU:HA	1:B:985:MET:HG3	1.79	0.62
1:A:1167:ARG:CZ	1:A:1176:ARG:HH12	2.12	0.62
1:A:325:ARG:NH2	2:A:1301:SO4:O3	2.32	0.62
1:A:696:VAL:HG22	1:A:769:ILE:HB	1.81	0.62
1:B:154:VAL:HB	1:B:239:ALA:HB3	1.81	0.62

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:666:ARG:HA	1:B:814:GLU:O	2.00	0.62
1:A:1064:ASP:OD1	1:A:1067:ARG:NH2	2.31	0.62
1:A:724:GLY:HA2	1:A:750:ILE:O	1.99	0.61
1:A:870:ARG:NH2	1:A:928:GLU:OE2	2.33	0.61
1:A:734:SER:OG	1:A:754:ARG:NH2	2.33	0.61
1:B:985:MET:O	1:B:988:ALA:N	2.33	0.61
1:B:857:ILE:HD13	1:B:941:LEU:HD13	1.83	0.61
1:A:1091:ILE:HD13	1:A:1096:PRO:HA	1.83	0.60
1:B:969:TYR:O	1:B:973:ALA:HB2	2.01	0.60
1:A:1120:THR:HB	1:A:1134:PRO:HG3	1.83	0.60
1:A:185:ARG:NH2	1:A:202:ALA:O	2.35	0.60
1:A:1073:ARG:NH2	1:A:1089:GLU:OE2	2.35	0.59
1:A:364:THR:HG23	1:A:367:GLU:H	1.67	0.59
1:B:72:ALA:O	1:B:76:THR:OG1	2.17	0.59
1:A:92:SER:HB2	1:A:138:ARG:HD2	1.85	0.58
1:A:335:GLU:OE2	1:A:845:ARG:NH2	2.36	0.58
1:B:752:THR:OG1	1:B:753:HIS:N	2.34	0.58
1:A:836:ASP:HB3	1:A:839:GLN:HG3	1.87	0.57
1:A:427:VAL:HG11	1:A:438:VAL:HG11	1.86	0.57
1:B:611:VAL:HG21	1:B:1028:TYR:HB3	1.86	0.57
1:B:911:THR:HG22	1:B:941:LEU:HD21	1.87	0.57
1:B:862:ARG:HG3	1:B:863:THR:HG23	1.87	0.56
1:B:969:TYR:O	1:B:973:ALA:CB	2.53	0.56
1:A:65:VAL:HG21	1:A:140:LEU:HD22	1.87	0.56
1:A:935:LEU:HG	1:A:971:ARG:HG3	1.86	0.56
1:A:522:VAL:HG22	1:A:528:ILE:HG12	1.87	0.56
1:B:442:LEU:HD12	1:B:449:ALA:HB2	1.88	0.56
1:A:856:TYR:O	1:A:908:VAL:HA	2.06	0.56
1:B:1165:LEU:O	1:B:1176:ARG:NH1	2.39	0.56
1:A:168:VAL:HG23	1:A:188:PHE:CE2	2.41	0.55
1:A:685:PHE:HD2	1:A:715:MET:HG2	1.71	0.55
1:A:911:THR:HG22	1:A:941:LEU:HD21	1.89	0.55
1:A:174:LEU:HD22	1:A:241:PRO:HG3	1.88	0.55
1:A:255:ALA:O	1:A:259:ALA:CB	2.54	0.55
1:B:47:SER:HB3	1:B:472:SER:HB2	1.88	0.55
1:B:681:VAL:HG21	1:B:714:ARG:HD3	1.87	0.55
1:A:13:ILE:HG13	1:A:82:VAL:HG12	1.87	0.55
1:A:168:VAL:HG23	1:A:188:PHE:HE2	1.70	0.55
1:A:727:ARG:NH2	1:A:893:GLU:OE1	2.40	0.55
1:B:856:TYR:O	1:B:908:VAL:HA	2.06	0.55
1:B:208:VAL:HG12	1:B:218:MET:HG2	1.88	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:335:GLU:OE2	1:B:845:ARG:NH2	2.40	0.54
1:B:726:SER:O	1:B:729:THR:OG1	2.24	0.54
1:A:524:ASP:OD2	1:A:1019:ARG:NH2	2.41	0.54
1:A:163:GLU:O	1:A:167:VAL:HG23	2.08	0.54
1:B:701:THR:HG21	1:B:727:ARG:H	1.71	0.54
1:A:26:LEU:HD21	1:A:395:ILE:HD11	1.90	0.53
1:B:857:ILE:HD12	1:B:925:LEU:HD11	1.90	0.53
1:B:94:GLU:HG2	1:B:951:ARG:HE	1.72	0.53
1:B:1052:VAL:HG13	1:B:1179:ASP:HA	1.91	0.53
1:B:811:GLY:HA3	1:B:1029:ARG:HE	1.74	0.53
1:B:624:ALA:HB2	1:B:661:PRO:HA	1.91	0.52
1:A:603:VAL:HG21	1:A:1021:VAL:HA	1.92	0.52
1:B:571:ASP:N	1:B:571:ASP:OD1	2.39	0.52
1:B:857:ILE:HD11	1:B:914:GLU:HG3	1.92	0.52
1:A:146:PRO:HD2	1:A:299:ILE:HD13	1.91	0.52
1:B:1191:LEU:HD13	1:B:1200:ASP:HB2	1.92	0.52
1:B:137:THR:HG22	1:B:322:VAL:HG13	1.92	0.52
1:B:712:THR:HG22	1:B:722:VAL:HG21	1.91	0.52
1:B:453:GLU:HG3	1:B:456:THR:HG21	1.90	0.52
1:A:49:ARG:HE	1:A:317:CYS:HB3	1.74	0.51
1:A:935:LEU:H	1:A:935:LEU:HD12	1.75	0.51
1:A:1087:VAL:HG13	1:A:1097:LEU:HD11	1.92	0.51
1:A:730:ASP:OD1	1:A:730:ASP:N	2.42	0.51
1:B:17:VAL:HG21	1:B:82:VAL:HG11	1.90	0.51
1:B:665:ASP:OD1	1:B:813:ARG:NE	2.44	0.51
1:A:262:HIS:NE2	1:A:349:ASP:O	2.43	0.51
1:B:990:LYS:HA	1:B:993:GLU:HG3	1.92	0.51
1:A:217:GLU:OE1	1:A:219:ARG:NH2	2.44	0.51
1:A:342:SER:O	1:A:346:VAL:HG23	2.10	0.51
1:B:180:ASP:O	1:B:191:ARG:NH1	2.22	0.51
1:B:63:LEU:HD22	1:B:308:LEU:HD11	1.93	0.51
1:B:65:VAL:HG21	1:B:140:LEU:HD22	1.91	0.51
1:A:651:ILE:HG22	1:A:655:LYS:HE3	1.92	0.50
1:B:880:VAL:HG22	1:B:906:ILE:HB	1.93	0.50
1:B:156:LEU:HB2	1:B:237:VAL:HG22	1.94	0.50
1:A:1147:ARG:NH1	2:A:1303:SO4:O1	2.38	0.50
1:A:100:ARG:NH2	1:A:340:SER:O	2.44	0.50
1:B:828:VAL:HG12	1:B:946:GLY:HA3	1.94	0.50
1:A:1133:SER:HB3	1:A:1134:PRO:HD3	1.93	0.50
1:B:194:ILE:HG12	1:B:209:GLU:HG3	1.92	0.50
1:A:803:ARG:HA	1:A:806:GLU:HB3	1.94	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:76:THR:HG21	1:A:89:LEU:HB2	1.94	0.49
1:A:1156:ALA:N	2:A:1302:SO4:O4	2.44	0.49
1:A:594:ASN:O	1:A:598:LYS:HG2	2.13	0.49
1:A:79:LEU:HD13	1:A:132:VAL:HG11	1.95	0.49
1:A:929:ARG:NH2	5:A:1403:HOH:O	2.46	0.48
1:B:21:LEU:HD23	1:B:26:LEU:HD13	1.95	0.48
1:B:52:VAL:O	1:B:55:ALA:HB3	2.12	0.48
1:A:1120:THR:H	1:A:1134:PRO:CG	2.26	0.48
1:A:565:VAL:HG21	1:A:573:LEU:HD11	1.95	0.48
1:A:991:ASP:HA	1:A:994:ILE:HG12	1.95	0.48
1:B:26:LEU:O	1:B:30:ILE:HG12	2.14	0.48
1:A:928:GLU:HG2	1:A:929:ARG:HG2	1.95	0.48
1:B:1054:ALA:HB1	1:B:1104:LEU:HA	1.95	0.48
1:B:330:ILE:HD11	1:B:364:THR:HA	1.96	0.48
1:A:415:LEU:HD22	1:A:425:ALA:HB1	1.96	0.47
1:B:13:ILE:HG13	1:B:82:VAL:HG12	1.96	0.47
1:B:832:VAL:HA	1:B:956:PHE:O	2.14	0.47
1:A:325:ARG:O	1:A:329:LEU:HD22	2.13	0.47
1:A:681:VAL:HG12	1:A:711:PHE:CE1	2.50	0.47
1:B:287:ASP:OD1	1:B:288:GLY:N	2.47	0.47
1:B:847:GLU:OE1	1:B:924:THR:OG1	2.29	0.47
1:B:49:ARG:HE	1:B:317:CYS:HB3	1.77	0.47
1:B:396:ARG:HG3	1:B:475:VAL:HB	1.97	0.47
1:A:1061:ILE:O	1:A:1067:ARG:NH1	2.48	0.47
1:A:802:PRO:HD3	1:A:993:GLU:HB3	1.96	0.47
1:B:419:VAL:HG11	1:B:463:VAL:HG22	1.97	0.47
1:A:647:GLN:O	1:A:651:ILE:HG13	2.15	0.47
1:B:1129:THR:HA	1:B:1164:PRO:HA	1.97	0.47
1:B:882:VAL:HA	1:B:908:VAL:O	2.15	0.47
1:A:970:ASP:O	1:A:974:THR:HG23	2.14	0.47
1:A:161:GLU:HA	1:A:214:GLU:HB3	1.97	0.47
1:A:415:LEU:HD21	1:A:484:ILE:HD12	1.97	0.47
4:A:1305:ADP:C8	4:A:1305:ADP:H5'2	2.50	0.46
1:A:931:ASP:OD1	1:A:931:ASP:N	2.46	0.46
1:B:1045:ASP:N	1:B:1045:ASP:OD1	2.48	0.46
1:B:198:PHE:CD2	1:B:205:PRO:HG3	2.51	0.46
1:B:1137:LEU:HB3	1:B:1141:ALA:HB3	1.97	0.46
1:B:141:LEU:HD11	1:B:326:ALA:HB1	1.97	0.46
1:A:822:PRO:HB2	4:A:1305:ADP:N1	2.31	0.46
1:A:888:ASN:N	1:A:888:ASN:OD1	2.49	0.46
1:B:322:VAL:HG12	1:B:365:PHE:HE2	1.81	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:406:HIS:CE1	1:B:396:ARG:HD2	2.51	0.46
1:B:407:ASN:O	1:B:411:ILE:HG12	2.15	0.46
1:B:696:VAL:HG22	1:B:769:ILE:HB	1.97	0.46
1:A:636:GLU:HG2	1:A:682:ARG:HD3	1.98	0.46
1:B:290:GLU:O	1:B:293:LEU:HB2	2.16	0.46
1:A:476:LEU:HD13	1:A:481:LEU:HD23	1.98	0.45
1:A:980:GLU:HA	1:A:985:MET:SD	2.56	0.45
1:B:439:VAL:HG22	1:B:449:ALA:HB1	1.98	0.45
1:B:453:GLU:O	1:B:456:THR:OG1	2.24	0.45
1:A:28:ASP:OD1	1:A:31:ARG:NH2	2.49	0.45
1:A:584:LEU:HD23	1:A:584:LEU:HA	1.79	0.45
1:B:369:ARG:HG3	1:B:379:TRP:CG	2.51	0.45
1:B:788:ARG:HG2	1:B:793:VAL:HG11	1.97	0.45
1:B:1130:VAL:HB	1:B:1165:LEU:HD21	1.98	0.45
1:A:23:ASP:HB2	1:A:479:ALA:HB2	1.99	0.45
1:A:387:ASP:OD1	1:A:388:GLU:N	2.48	0.45
1:A:666:ARG:HA	1:A:814:GLU:O	2.16	0.45
1:A:86:SER:HB3	1:A:131:ARG:HG3	1.97	0.45
1:B:528:ILE:HD12	1:B:583:SER:HB2	1.97	0.45
1:B:937:GLN:O	1:B:941:LEU:HG	2.17	0.45
1:B:1133:SER:HB3	1:B:1134:PRO:HD3	1.98	0.45
1:B:305:THR:HG21	1:B:372:ALA:HB2	1.98	0.45
1:A:402:ARG:NH2	1:B:397:SER:OG	2.30	0.45
1:A:1114:CYS:HB3	1:A:1119:ILE:HB	1.97	0.45
1:B:440:GLU:O	1:B:444:GLU:HB2	2.17	0.45
1:B:704:ALA:O	1:B:708:LEU:CB	2.64	0.45
1:A:326:ALA:HA	1:A:329:LEU:HD23	1.99	0.44
1:B:342:SER:O	1:B:346:VAL:HG23	2.17	0.44
1:B:42:LEU:HD22	1:B:393:LEU:HD11	1.99	0.44
1:B:770:VAL:HB	1:B:795:THR:HG22	1.98	0.44
1:B:935:LEU:H	1:B:935:LEU:HD12	1.82	0.44
1:A:1149:TYR:HE1	1:A:1192:VAL:HG21	1.82	0.44
1:A:535:THR:OG1	1:A:536:GLU:N	2.50	0.44
1:A:643:GLU:CD	1:A:643:GLU:H	2.20	0.44
1:A:743:ASP:OD1	1:A:743:ASP:N	2.50	0.44
1:A:677:THR:O	1:A:681:VAL:HG13	2.17	0.44
1:A:904:TYR:CE1	1:B:877:GLU:HB3	2.48	0.44
1:A:40:LEU:HB3	1:A:380:TRP:CG	2.52	0.44
1:B:1087:VAL:HG13	1:B:1097:LEU:HD11	1.99	0.44
1:B:285:PRO:HB3	1:B:289:MET:HE1	1.99	0.44
1:A:937:GLN:O	1:A:941:LEU:HG	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:100:ARG:NH2	1:B:271:GLY:O	2.45	0.44
1:A:704:ALA:O	1:A:708:LEU:HB2	2.18	0.44
1:B:931:ASP:N	1:B:931:ASP:OD1	2.48	0.44
1:A:830:THR:OG1	1:A:942:ARG:NH1	2.50	0.43
1:B:1060:TYR:CZ	1:B:1094:TYR:HB2	2.53	0.43
1:B:364:THR:HG23	1:B:367:GLU:H	1.83	0.43
1:B:853:GLN:NE2	1:B:899:PHE:O	2.49	0.43
1:A:47:SER:HB3	1:A:472:SER:HB2	1.99	0.43
1:B:321:LYS:HD2	1:B:321:LYS:H	1.84	0.43
1:B:68:THR:OG1	1:B:71:GLU:OE1	2.22	0.43
1:A:888:ASN:ND2	1:A:890:GLU:OE1	2.46	0.43
1:B:1161:VAL:HG11	1:B:1193:LEU:HD11	2.00	0.43
1:B:866:GLU:O	1:B:870:ARG:HG3	2.18	0.43
1:B:1149:TYR:HE2	1:B:1192:VAL:HG21	1.83	0.43
1:B:596:LYS:HB3	1:B:596:LYS:HE2	1.78	0.43
1:B:666:ARG:NH2	1:B:816:SER:OG	2.51	0.43
1:A:17:VAL:HG21	1:A:82:VAL:HG11	2.00	0.43
1:B:1122:ILE:HG12	1:B:1132:LEU:HG	2.00	0.43
1:B:76:THR:HG21	1:B:89:LEU:HB2	2.00	0.43
1:A:272:THR:HG23	1:A:274:PRO:HD2	2.00	0.43
1:B:90:PHE:HB3	1:B:135:THR:HB	2.01	0.43
1:A:290:GLU:HA	1:A:293:LEU:HD13	2.00	0.43
1:B:643:GLU:N	1:B:643:GLU:OE1	2.40	0.43
1:A:722:VAL:HG12	1:A:748:VAL:CG1	2.49	0.43
1:B:276:MET:HG3	1:B:286:VAL:HG21	2.01	0.43
1:A:829:LEU:HD12	1:A:953:TYR:HE1	1.83	0.42
1:A:990:LYS:O	1:A:994:ILE:HG12	2.19	0.42
4:B:1302:ADP:H5'2	4:B:1302:ADP:H8	1.83	0.42
1:B:308:LEU:HB2	1:B:377:HIS:CE1	2.54	0.42
1:A:23:ASP:OD2	1:A:25:SER:HB3	2.19	0.42
1:B:152:GLU:HB3	1:B:170:ARG:HH11	1.84	0.42
1:A:165:GLU:HA	1:A:168:VAL:HG12	2.00	0.42
1:B:92:SER:HB2	1:B:138:ARG:HD2	2.01	0.42
1:A:536:GLU:HA	1:A:544:ARG:O	2.19	0.42
1:A:574:SER:HB2	1:A:1015:ASP:HB3	2.02	0.42
1:A:972:LEU:O	1:A:976:ALA:HB2	2.18	0.42
1:B:990:LYS:O	1:B:994:ILE:HG12	2.20	0.42
1:A:87:VAL:HA	1:A:132:VAL:O	2.20	0.42
1:B:126:LEU:HD23	1:B:238:VAL:HG21	2.00	0.42
1:A:626:GLY:O	1:A:686:LYS:NZ	2.50	0.42
1:B:725:LEU:HD11	1:B:754:ARG:HE	1.85	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1198:GLN:N	1:A:1198:GLN:OE1	2.51	0.42
1:A:154:VAL:HB	1:A:239:ALA:HB3	2.02	0.42
1:A:452:LEU:HB2	1:A:466:LEU:HD13	2.00	0.42
1:A:653:GLU:OE2	1:A:816:SER:OG	2.26	0.42
1:A:694:VAL:O	1:A:748:VAL:HA	2.20	0.42
1:B:106:GLU:HG2	1:B:106:GLU:H	1.69	0.42
1:A:538:VAL:HA	1:A:542:ALA:O	2.20	0.42
1:A:645:ILE:HG23	1:A:820:THR:HG21	2.02	0.42
1:B:788:ARG:HH21	1:B:795:THR:HG21	1.84	0.42
1:A:63:LEU:HD12	1:A:133:VAL:O	2.20	0.41
1:A:63:LEU:HD22	1:A:308:LEU:HD11	2.01	0.41
1:B:63:LEU:HD12	1:B:133:VAL:O	2.18	0.41
1:B:1180:LEU:HD23	1:B:1180:LEU:HA	1.75	0.41
1:B:674:TYR:CE1	1:B:821:PRO:HA	2.55	0.41
1:B:856:TYR:HB3	1:B:908:VAL:HG22	2.01	0.41
1:A:708:LEU:O	1:A:712:THR:HG23	2.20	0.41
1:A:756:LEU:HD12	1:A:780:HIS:HB3	2.02	0.41
1:B:56:LEU:HB2	1:B:62:LEU:HD13	2.01	0.41
1:B:94:GLU:HG2	1:B:951:ARG:NE	2.34	0.41
1:A:94:GLU:OE2	1:A:107:THR:OG1	2.30	0.41
1:A:224:ALA:HA	1:A:509:VAL:H	1.85	0.41
1:A:23:ASP:HA	1:A:24:PRO:HD3	1.96	0.41
1:B:1091:ILE:HA	1:B:1095:GLY:O	2.20	0.41
1:A:428:VAL:HG21	1:A:470:LEU:HD12	2.03	0.41
1:A:93:TRP:NE1	1:A:244:GLU:OE2	2.52	0.41
1:B:135:THR:OG1	1:B:136:THR:N	2.54	0.41
1:B:721:THR:OG1	1:B:747:ASP:N	2.54	0.41
1:B:1188:GLY:O	1:B:1192:VAL:HG22	2.21	0.41
1:B:248:THR:O	1:B:251:VAL:HG22	2.21	0.41
1:B:670:GLY:O	1:B:798:ALA:HA	2.21	0.41
1:A:141:LEU:O	1:A:363:VAL:HG22	2.21	0.41
1:B:195:LEU:O	1:B:208:VAL:HG22	2.21	0.41
1:A:770:VAL:HB	1:A:795:THR:HG22	2.02	0.40
1:A:452:LEU:HD12	1:A:466:LEU:HB2	2.03	0.40
1:A:841:ALA:O	1:A:845:ARG:HG3	2.22	0.40
1:B:859:ASN:ND2	5:B:1401:HOH:O	2.36	0.40
1:A:1093:ARG:NH1	1:A:1094:TYR:OH	2.54	0.40
1:A:273:VAL:HG21	1:A:346:VAL:CG1	2.52	0.40
1:A:427:VAL:HA	1:A:484:ILE:O	2.21	0.40
1:A:607:ALA:O	1:A:611:VAL:HG22	2.21	0.40
1:B:708:LEU:O	1:B:712:THR:HG23	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:964:LEU:HD23	1:A:964:LEU:H	1.87	0.40
1:B:822:PRO:HB2	4:B:1302:ADP:C2	2.56	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	1165/1235 (94%)	1101 (94%)	54 (5%)	10 (1%)	17	56
1	B	1156/1235 (94%)	1088 (94%)	64 (6%)	4 (0%)	41	75
All	All	2321/2470 (94%)	2189 (94%)	118 (5%)	14 (1%)	25	64

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	511	ASP
1	A	1134	PRO
1	B	1134	PRO
1	A	512	PRO
1	A	1053	ASP
1	A	1175	PRO
1	B	224	ALA
1	A	824	GLU
1	A	224	ALA
1	A	263	PRO
1	A	984	GLY
1	A	192	GLY
1	B	346	VAL
1	B	451	ILE

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	851/988 (86%)	832 (98%)	19 (2%)	52	78
1	B	842/988 (85%)	827 (98%)	15 (2%)	59	81
All	All	1693/1976 (86%)	1659 (98%)	34 (2%)	55	79

All (34) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	27	GLN
1	A	164	PHE
1	A	198	PHE
1	A	262	HIS
1	A	287	ASP
1	A	289	MET
1	A	306	ARG
1	A	318	ASP
1	A	402	ARG
1	A	618	GLN
1	A	690	ASP
1	A	762	TRP
1	A	942	ARG
1	A	995	ARG
1	A	1014	PHE
1	A	1029	ARG
1	A	1079	ASP
1	A	1115	ARG
1	A	1203	MET
1	B	198	PHE
1	B	289	MET
1	B	414	MET
1	B	671	ASP
1	B	690	ASP
1	B	714	ARG
1	B	743	ASP
1	B	762	TRP

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Mol	Chain	Res	Type
1	B	776	PHE
1	B	972	LEU
1	B	1019	ARG
1	B	1080	ASP
1	B	1176	ARG
1	B	1200	ASP
1	B	1203	MET

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

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

7 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
2	SO4	B	1301	-	4,4,4	0.15	0	6,6,6	0.05	0
2	SO4	A	1302	-	4,4,4	0.13	0	6,6,6	0.05	0
4	ADP	B	1302	-	24,29,29	0.96	1 (4%)	29,45,45	1.49	4 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	SO4	A	1301	-	4,4,4	0.14	0	6,6,6	0.05	0
4	ADP	A	1305	-	24,29,29	0.94	1 (4%)	29,45,45	1.47	4 (13%)
3	PO4	A	1304	-	4,4,4	0.92	0	6,6,6	0.46	0
2	SO4	A	1303	-	4,4,4	0.14	0	6,6,6	0.05	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	ADP	A	1305	-	-	6/12/32/32	0/3/3/3
4	ADP	B	1302	-	-	5/12/32/32	0/3/3/3

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	1302	ADP	C5-C4	2.54	1.47	1.40
4	A	1305	ADP	C5-C4	2.44	1.47	1.40

All (8) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	1305	ADP	PA-O3A-PB	-3.87	119.54	132.83
4	B	1302	ADP	PA-O3A-PB	-3.62	120.41	132.83
4	B	1302	ADP	C3'-C2'-C1'	3.49	106.23	100.98
4	A	1305	ADP	N3-C2-N1	-3.16	123.74	128.68
4	B	1302	ADP	N3-C2-N1	-3.09	123.84	128.68
4	A	1305	ADP	C3'-C2'-C1'	2.53	104.78	100.98
4	B	1302	ADP	C4-C5-N7	-2.43	106.86	109.40
4	A	1305	ADP	C4-C5-N7	-2.28	107.02	109.40

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	1302	ADP	C5'-O5'-PA-O1A
4	B	1302	ADP	C5'-O5'-PA-O2A
4	A	1305	ADP	PB-O3A-PA-O5'
4	A	1305	ADP	C5'-O5'-PA-O3A

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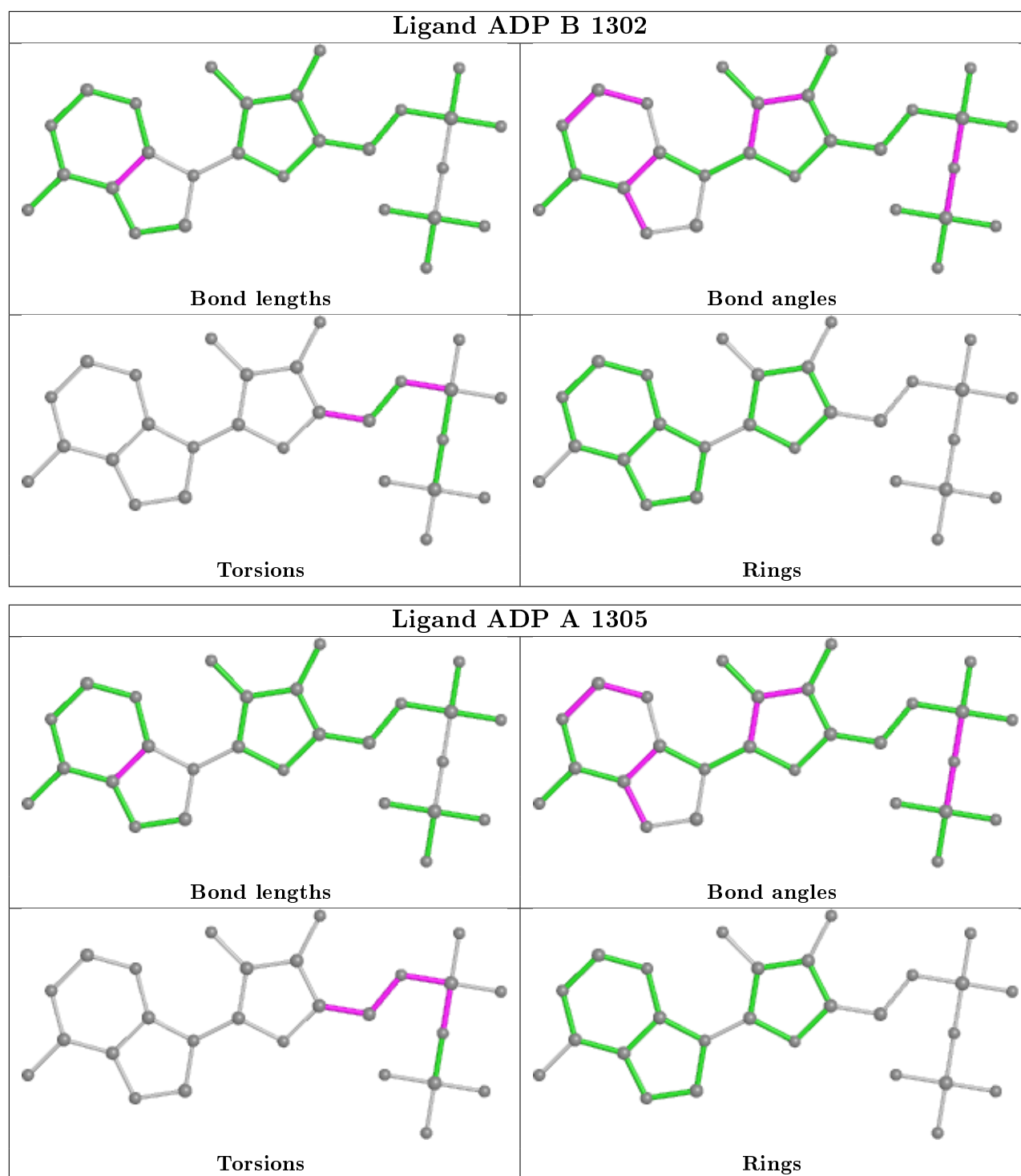
Mol	Chain	Res	Type	Atoms
4	A	1305	ADP	C3'-C4'-C5'-O5'
4	B	1302	ADP	C3'-C4'-C5'-O5'
4	A	1305	ADP	O4'-C4'-C5'-O5'
4	B	1302	ADP	O4'-C4'-C5'-O5'
4	A	1305	ADP	C4'-C5'-O5'-PA
4	A	1305	ADP	C5'-O5'-PA-O1A
4	B	1302	ADP	C5'-O5'-PA-O3A

There are no ring outliers.

5 monomers are involved in 8 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1302	SO4	1	0
4	B	1302	ADP	2	0
2	A	1301	SO4	1	0
4	A	1305	ADP	3	0
2	A	1303	SO4	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

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	1175/1235 (95%)	-0.37	12 (1%) 82 77	31, 64, 104, 132	0
1	B	1168/1235 (94%)	-0.30	15 (1%) 77 71	32, 71, 111, 135	0
All	All	2343/2470 (94%)	-0.34	27 (1%) 79 73	31, 67, 108, 135	0

All (27) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	8	SER	3.8
1	A	8	SER	3.5
1	B	507	ARG	3.2
1	A	622	GLY	2.9
1	A	580	GLU	2.9
1	B	561	ASP	2.8
1	B	404	SER	2.8
1	B	885	GLY	2.8
1	B	506	ARG	2.8
1	B	268	THR	2.7
1	A	504	ALA	2.7
1	A	541	GLY	2.7
1	A	621	PRO	2.5
1	B	822	PRO	2.5
1	B	821	PRO	2.5
1	A	964	LEU	2.5
1	A	1174	ALA	2.5
1	A	1031	ALA	2.4
1	B	505	LYS	2.4
1	A	822	PRO	2.2
1	A	982	GLY	2.2
1	B	406	HIS	2.1
1	B	282	GLU	2.1
1	B	884	HIS	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	1028	TYR	2.0
1	B	1033	ASP	2.0
1	B	970	ASP	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](#)

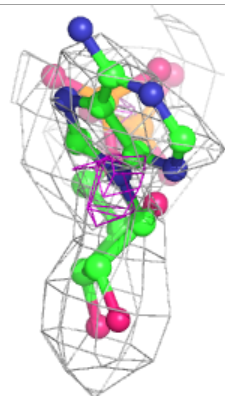
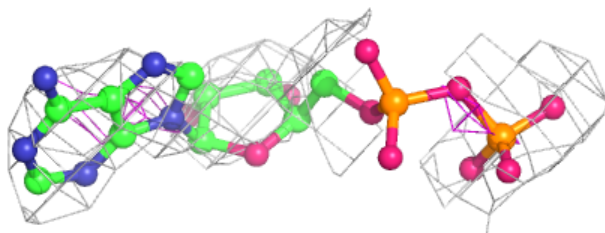
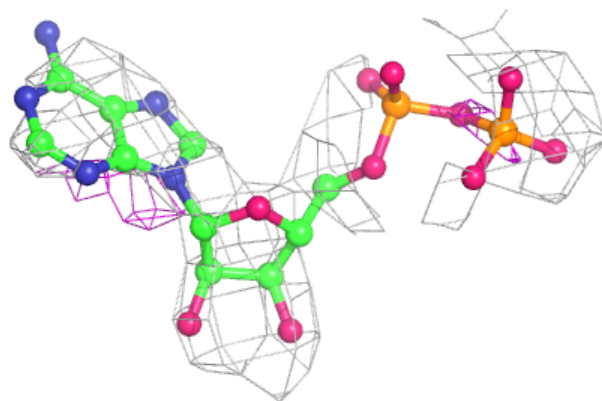
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
4	ADP	B	1302	27/27	0.64	0.40	81,103,115,136	0
4	ADP	A	1305	27/27	0.73	0.37	71,95,107,111	0
3	PO4	A	1304	5/5	0.81	0.40	87,94,97,105	0
2	SO4	A	1303	5/5	0.93	0.16	86,89,102,111	0
2	SO4	B	1301	5/5	0.94	0.17	48,56,63,66	0
2	SO4	A	1301	5/5	0.97	0.13	46,48,54,56	0
2	SO4	A	1302	5/5	0.97	0.12	52,64,65,68	0

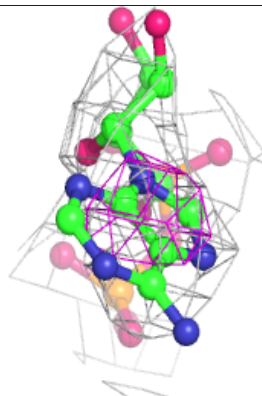
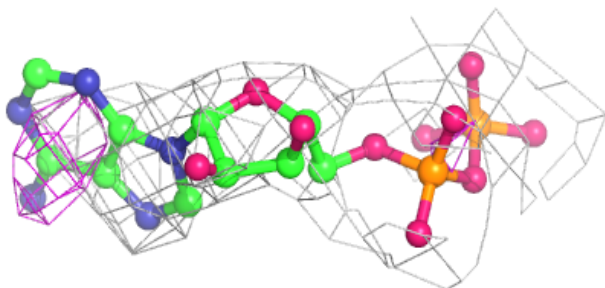
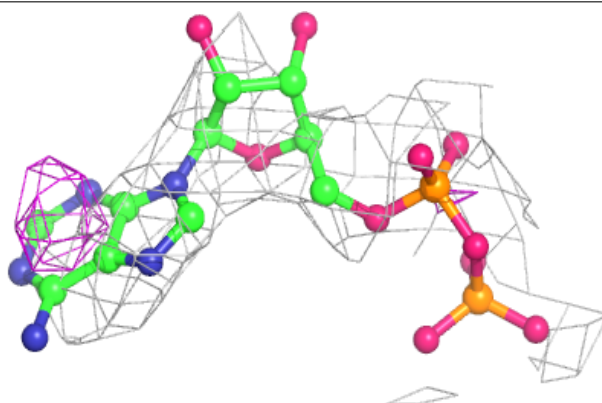
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

Electron density around ADP B 1302:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around ADP A 1305:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



6.5 Other polymers [i](#)

There are no such residues in this entry.