Full wwPDB X-ray Structure Validation Report (i)

Nov 2, 2023 – 02:57 AM EDT

PDB ID	:	3IKJ
Title	:	Structural characterization for the nucleotide binding ability of subunit A mutant S238A of the A1AO ATP synthase
Authors	:	Kumar, A.; Manimekali, M.S.S.; Balakrishna, A.M.; Jeyakanthan, J.; Gruber, G.
Deposited on	:	2009-08-06
Resolution	:	2.40 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 types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

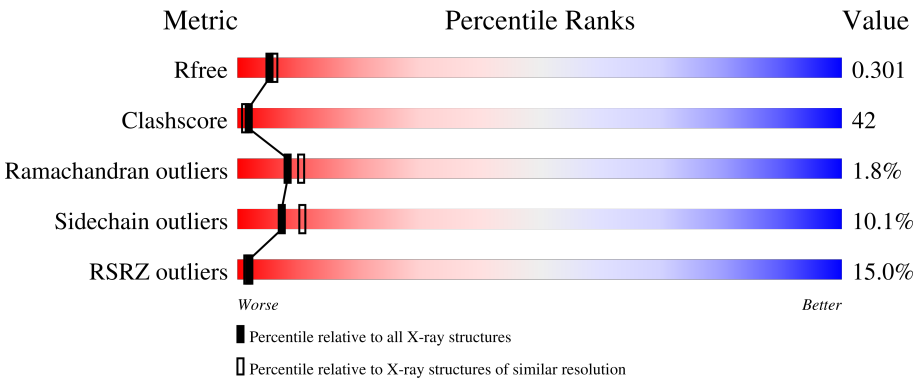
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

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 2.40 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	3907 (2.40-2.40)
Clashscore	141614	4398 (2.40-2.40)
Ramachandran outliers	138981	4318 (2.40-2.40)
Sidechain outliers	138945	4319 (2.40-2.40)
RSRZ outliers	127900	3811 (2.40-2.40)

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 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% 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	588	13% 45% 34% 7% 13%

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
2	MPD	A	589	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	MPD	A	591	-	-	X	-

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 4353 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.

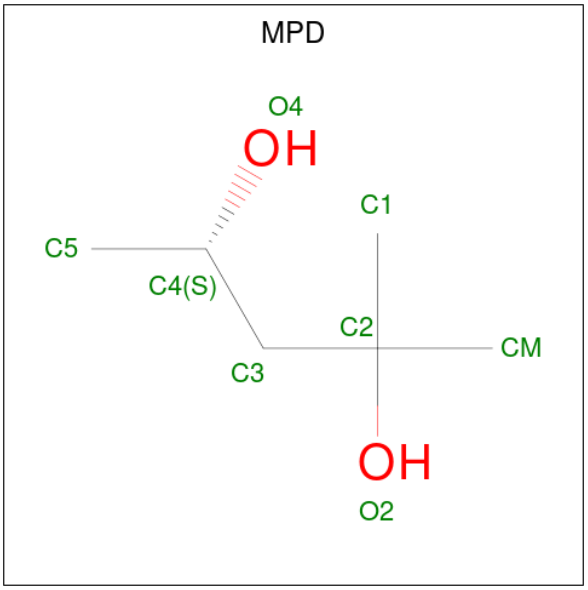
- Molecule 1 is a protein called V-type ATP synthase alpha chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	513	4055	2595	693	751	16	0	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	79	ARG	GLY	engineered mutation	UNP O57728
A	238	ALA	SER	engineered mutation	UNP O57728

- Molecule 2 is (4S)-2-METHYL-2,4-PENTANEDIOL (three-letter code: MPD) (formula: $C_6H_{14}O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	8	6	2	0	0
2	A	1	8	6	2	0	0

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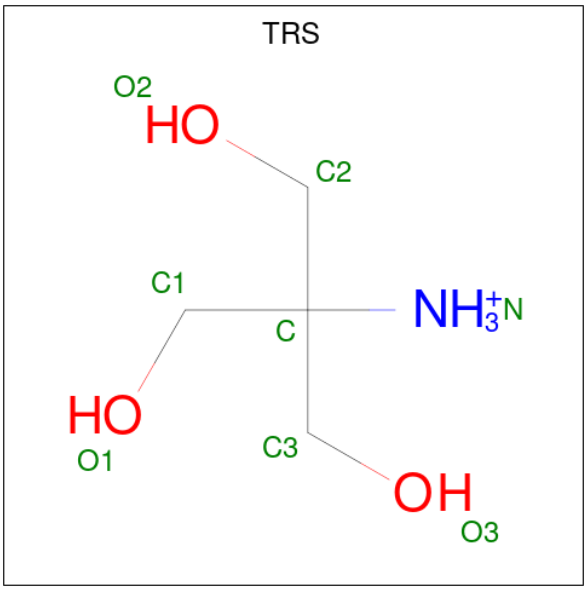
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
2	A	1	8	6	2	0	0

- Molecule 3 is 2-AMINO-2-HYDROXYMETHYL-PROPANE-1,3-DIOL (three-letter code: TRS) (formula: $C_4H_{12}NO_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	8	4	1	3	0	0

- Molecule 4 is water.

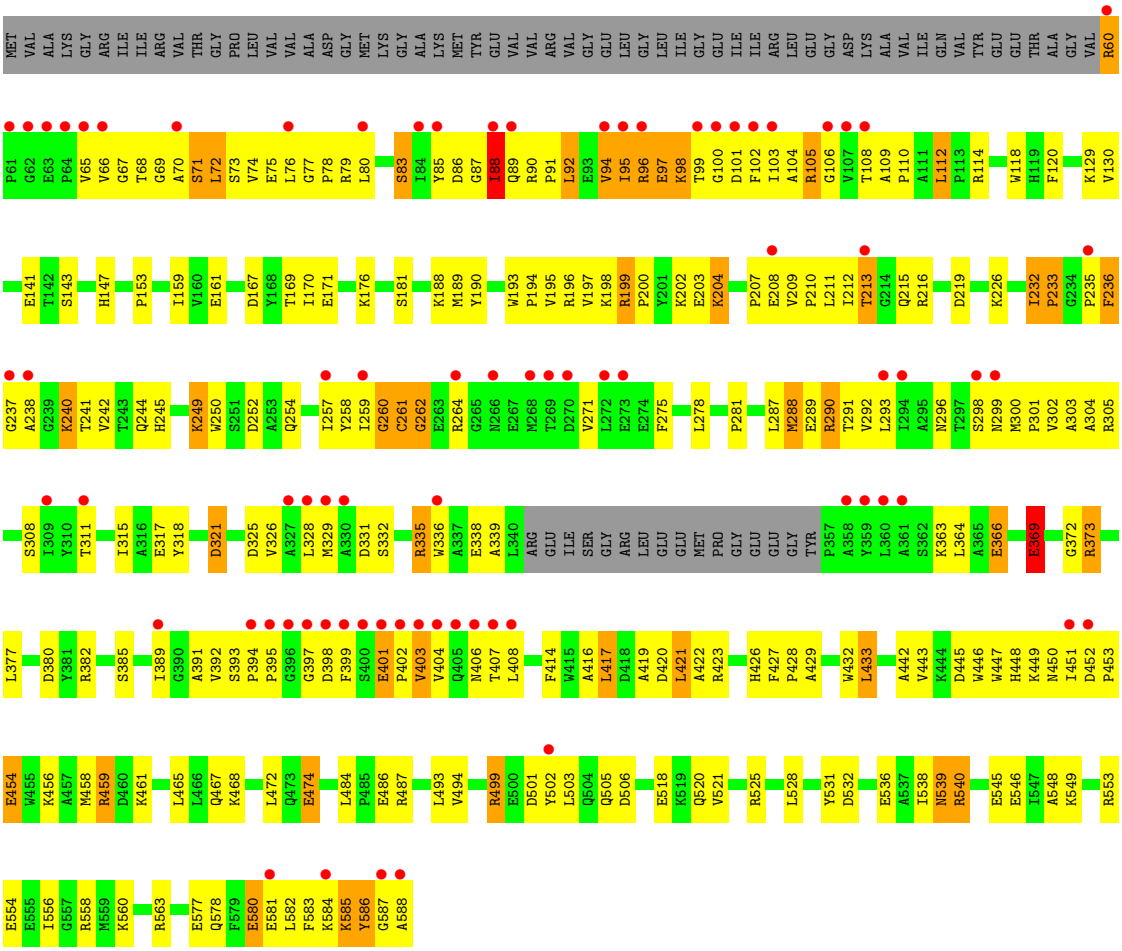
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	266	266	266	0	0

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 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: V-type ATP synthase alpha chain

Chain A: 13% 45% 34% 7% 13%



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α, β, γ	128.21Å 128.21Å 105.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.84 – 2.40 29.84 – 2.40	Depositor EDS
% Data completeness (in resolution range)	99.9 (29.84-2.40) 99.9 (29.84-2.40)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$< I/\sigma(I) >$ [1]	4.11 (at 2.39Å)	Xtriage
Refinement program	REFMAC 5.5.0072	Depositor
R, R_{free}	0.239 , 0.304 0.251 , 0.301	Depositor DCC
R_{free} test set	1752 reflections (5.02%)	wwPDB-VP
Wilson B-factor (Å^2)	50.8	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.31 , 62.2	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	4353	wwPDB-VP
Average B, all atoms (Å^2)	56.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.19% of the height of the origin peak. No significant pseudotranslation is detected.*

[1] Intensities estimated from amplitudes.

[2] Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 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: TRS, MPD

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	0.96	3/4144 (0.1%)	0.99	20/5614 (0.4%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand.A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	369	GLU	CG-CD	5.81	1.60	1.51
1	A	518	GLU	CG-CD	5.54	1.60	1.51
1	A	454	GLU	CG-CD	5.30	1.59	1.51

All (20) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	373	ARG	NE-CZ-NH2	-13.09	113.76	120.30
1	A	499	ARG	NE-CZ-NH2	-11.60	114.50	120.30
1	A	199	ARG	NE-CZ-NH1	-10.44	115.08	120.30
1	A	465	LEU	CB-CG-CD2	-9.20	95.36	111.00
1	A	199	ARG	NE-CZ-NH2	9.01	124.80	120.30
1	A	373	ARG	NE-CZ-NH1	8.03	124.31	120.30
1	A	321	ASP	CB-CG-OD1	7.20	124.78	118.30
1	A	506	ASP	CB-CG-OD1	6.72	124.34	118.30
1	A	499	ARG	NE-CZ-NH1	6.63	123.62	120.30
1	A	459	ARG	CG-CD-NE	6.59	125.63	111.80

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	292	VAL	CB-CA-C	-6.07	99.87	111.40
1	A	484	LEU	CA-CB-CG	5.92	128.91	115.30
1	A	459	ARG	NE-CZ-NH2	5.89	123.25	120.30
1	A	459	ARG	N-CA-CB	5.74	120.93	110.60
1	A	112	LEU	CA-CB-CG	5.71	128.44	115.30
1	A	373	ARG	CG-CD-NE	-5.43	100.40	111.80
1	A	585	LYS	N-CA-C	-5.23	96.88	111.00
1	A	433	LEU	CA-CB-CG	5.15	127.15	115.30
1	A	335	ARG	NE-CZ-NH2	-5.11	117.75	120.30
1	A	459	ARG	CA-CB-CG	5.00	124.40	113.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	260	GLY	Peptide

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	4055	0	4112	345	0
2	A	24	0	42	27	0
3	A	8	0	12	1	0
4	A	266	0	0	43	0
All	All	4353	0	4166	350	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 42.

All (350) 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:261:CYS:SG	1:A:332:SER:HB2	1.41	1.56
1:A:233:PRO:HG3	1:A:414:PHE:CE1	1.55	1.38

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:HIS:NE2	2:A:589:MPD:H52	1.41	1.33
1:A:233:PRO:CG	1:A:414:PHE:CE1	2.15	1.27
1:A:97:GLU:HB2	1:A:105:ARG:O	1.27	1.25
1:A:261:CYS:SG	1:A:332:SER:CB	2.31	1.18
2:A:591:MPD:H53	4:A:690:HOH:O	1.39	1.18
1:A:83:SER:OG	1:A:85:TYR:CE2	2.00	1.15
1:A:79:ARG:HG2	4:A:818:HOH:O	0.97	1.14
1:A:79:ARG:HB2	1:A:118:TRP:CZ2	1.84	1.13
1:A:258:TYR:CE2	1:A:260:GLY:HA3	1.83	1.13
1:A:97:GLU:CB	1:A:105:ARG:O	1.98	1.12
1:A:96:ARG:HG3	1:A:301:PRO:HD2	1.22	1.11
1:A:90:ARG:NH2	1:A:96:ARG:HA	1.67	1.10
1:A:99:THR:HG22	1:A:100:GLY:H	1.15	1.10
1:A:88:ILE:HG22	1:A:89:GLN:H	1.05	1.08
1:A:245:HIS:CD2	1:A:249:LYS:HD3	1.89	1.08
1:A:233:PRO:HB3	1:A:392:VAL:HG23	1.37	1.07
1:A:79:ARG:O	1:A:79:ARG:HD3	1.54	1.06
1:A:233:PRO:CD	1:A:414:PHE:CE1	2.40	1.05
1:A:96:ARG:CG	1:A:301:PRO:HD2	1.87	1.04
1:A:389:ILE:HB	4:A:838:HOH:O	1.55	1.03
1:A:233:PRO:CG	1:A:414:PHE:CZ	2.42	1.02
1:A:88:ILE:CG2	1:A:89:GLN:H	1.72	1.02
1:A:90:ARG:HH12	1:A:96:ARG:HE	1.07	1.01
1:A:79:ARG:HB2	1:A:118:TRP:HZ2	1.13	1.01
1:A:99:THR:CG2	1:A:100:GLY:H	1.73	1.01
1:A:233:PRO:CG	1:A:414:PHE:HE1	1.59	1.00
1:A:233:PRO:HG3	1:A:414:PHE:HE1	0.90	1.00
1:A:80:LEU:O	1:A:83:SER:HB3	1.63	0.98
1:A:401:GLU:HG2	1:A:402:PRO:HD3	1.46	0.98
1:A:445:ASP:HB2	4:A:689:HOH:O	1.62	0.97
1:A:448:HIS:HA	1:A:452:ASP:O	1.65	0.97
1:A:245:HIS:CD2	2:A:589:MPD:H31	2.00	0.97
1:A:88:ILE:HG22	1:A:89:GLN:N	1.67	0.96
1:A:258:TYR:CE2	1:A:260:GLY:CA	2.48	0.96
1:A:75:GLU:H	1:A:89:GLN:HE22	1.03	0.94
1:A:245:HIS:CD2	2:A:589:MPD:H52	2.02	0.93
1:A:208:GLU:O	1:A:443:VAL:HG12	1.68	0.93
1:A:96:ARG:HG3	1:A:301:PRO:CD	1.99	0.93
1:A:261:CYS:HG	1:A:332:SER:HB2	1.28	0.92
1:A:213:THR:HG22	1:A:215:GLN:H	1.35	0.91
1:A:249:LYS:HE2	2:A:589:MPD:H53	1.50	0.91

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:216:ARG:H	1:A:505:GLN:HE22	1.17	0.90
1:A:232:ILE:HD13	1:A:391:ALA:HA	1.51	0.90
1:A:90:ARG:HH22	1:A:96:ARG:HA	1.33	0.90
1:A:546:GLU:HG3	1:A:582:LEU:HD11	1.53	0.89
1:A:112:LEU:HG	4:A:831:HOH:O	1.72	0.88
1:A:90:ARG:HH12	1:A:96:ARG:NE	1.71	0.88
1:A:258:TYR:HE2	1:A:260:GLY:HA3	1.35	0.87
1:A:120:PHE:HB2	4:A:756:HOH:O	1.73	0.86
1:A:96:ARG:HB2	1:A:301:PRO:HG2	1.56	0.86
1:A:233:PRO:HG3	1:A:414:PHE:CZ	2.08	0.86
1:A:233:PRO:HD3	1:A:414:PHE:CE1	2.10	0.85
1:A:102:PHE:C	1:A:103:ILE:HD12	1.98	0.84
1:A:458:MET:CE	1:A:525:ARG:HG2	2.07	0.84
1:A:216:ARG:HD2	1:A:502:TYR:CE1	2.13	0.84
1:A:233:PRO:HG2	1:A:414:PHE:CZ	2.12	0.84
1:A:96:ARG:CB	1:A:301:PRO:HD2	2.09	0.83
1:A:102:PHE:HB2	1:A:103:ILE:HD12	1.61	0.83
1:A:99:THR:HG22	1:A:100:GLY:N	1.93	0.83
1:A:91:PRO:O	1:A:94:VAL:HG22	1.78	0.83
1:A:108:THR:HG22	1:A:109:ALA:H	1.42	0.82
1:A:245:HIS:HD2	1:A:249:LYS:HD3	1.42	0.82
1:A:92:LEU:HD13	4:A:726:HOH:O	1.78	0.82
1:A:249:LYS:NZ	1:A:281:PRO:HG3	1.96	0.81
1:A:467:GLN:HG3	4:A:701:HOH:O	1.80	0.81
1:A:238:ALA:HB3	1:A:422:ALA:HB1	1.63	0.80
1:A:245:HIS:HD2	2:A:589:MPD:H31	1.43	0.80
1:A:96:ARG:HB2	1:A:301:PRO:HD2	1.64	0.80
1:A:249:LYS:CE	2:A:589:MPD:H53	2.12	0.79
1:A:216:ARG:HB3	1:A:502:TYR:CE2	2.19	0.78
1:A:86:ASP:O	1:A:88:ILE:O	2.00	0.78
1:A:79:ARG:O	1:A:79:ARG:CD	2.30	0.77
1:A:96:ARG:HB2	1:A:301:PRO:CG	2.15	0.77
1:A:74:VAL:HG23	1:A:195:VAL:CG1	2.15	0.76
1:A:245:HIS:NE2	2:A:589:MPD:C5	2.37	0.76
1:A:199:ARG:NH1	1:A:321:ASP:OD2	2.18	0.76
1:A:339:ALA:HA	4:A:824:HOH:O	1.84	0.76
1:A:580:GLU:O	1:A:584:LYS:HG3	1.86	0.76
1:A:96:ARG:HD3	1:A:300:MET:HG2	1.68	0.76
1:A:103:ILE:HG22	1:A:104:ALA:N	2.01	0.76
1:A:153:PRO:HG2	1:A:193:TRP:CZ3	2.21	0.76
1:A:88:ILE:CG2	1:A:89:GLN:N	2.35	0.75

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:78:PRO:O	4:A:831:HOH:O	2.03	0.75
1:A:454:GLU:HG3	4:A:830:HOH:O	1.86	0.75
1:A:261:CYS:HB3	1:A:296:ASN:H	1.51	0.74
2:A:589:MPD:HM1	2:A:589:MPD:O4	1.88	0.74
1:A:69:GLY:HA2	1:A:103:ILE:HA	1.71	0.73
1:A:108:THR:HG22	1:A:109:ALA:N	2.03	0.73
1:A:109:ALA:HB1	1:A:110:PRO:HD2	1.69	0.73
1:A:235:PRO:HG3	1:A:419:ALA:HB2	1.69	0.73
1:A:90:ARG:CZ	1:A:95:ILE:O	2.36	0.72
1:A:216:ARG:H	1:A:505:GLN:NE2	1.86	0.72
1:A:87:GLY:HA2	1:A:308:SER:HB3	1.70	0.72
1:A:96:ARG:HB2	1:A:301:PRO:CD	2.19	0.72
1:A:103:ILE:CG2	1:A:104:ALA:N	2.52	0.72
1:A:448:HIS:HE1	1:A:456:LYS:H	1.35	0.72
1:A:249:LYS:CD	2:A:589:MPD:H53	2.19	0.72
1:A:97:GLU:HB3	1:A:106:GLY:HA2	1.72	0.71
1:A:254:GLN:HE22	1:A:325:ASP:H	1.37	0.71
1:A:447:TRP:CZ3	1:A:451:ILE:HG21	2.25	0.71
1:A:311:THR:O	1:A:315:ILE:HG13	1.91	0.71
1:A:264:ARG:HD2	4:A:815:HOH:O	1.91	0.71
1:A:258:TYR:CD2	1:A:260:GLY:HA3	2.26	0.70
1:A:402:PRO:O	1:A:403:VAL:HG12	1.92	0.70
1:A:216:ARG:HD2	1:A:502:TYR:CD1	2.27	0.69
1:A:233:PRO:HG2	1:A:414:PHE:HZ	1.56	0.69
1:A:90:ARG:NH1	1:A:96:ARG:HE	1.87	0.69
1:A:366:GLU:N	1:A:366:GLU:OE2	2.25	0.69
1:A:452:ASP:OD1	1:A:453:PRO:N	2.25	0.69
1:A:75:GLU:N	1:A:89:GLN:HE22	1.86	0.69
1:A:69:GLY:HA2	1:A:103:ILE:HG13	1.75	0.69
1:A:143:SER:OG	1:A:289:GLU:OE2	2.11	0.68
1:A:521:VAL:O	1:A:525:ARG:HG3	1.94	0.68
1:A:90:ARG:HB2	1:A:94:VAL:HG21	1.76	0.68
1:A:563:ARG:HD3	4:A:762:HOH:O	1.94	0.68
1:A:259:ILE:O	1:A:261:CYS:SG	2.52	0.67
1:A:401:GLU:CG	1:A:402:PRO:HD3	2.24	0.67
1:A:91:PRO:O	1:A:94:VAL:CG2	2.43	0.67
1:A:211:LEU:HD12	1:A:252:ASP:OD1	1.94	0.66
1:A:427:PHE:HA	1:A:428:PRO:C	2.16	0.66
1:A:329:MET:HG2	4:A:838:HOH:O	1.95	0.66
1:A:83:SER:OG	1:A:85:TYR:CZ	2.37	0.66
1:A:461:LYS:HZ3	2:A:591:MPD:HM1	1.58	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:536:GLU:O	1:A:540:ARG:HG3	1.94	0.65
1:A:369:GLU:HB3	4:A:725:HOH:O	1.96	0.65
1:A:245:HIS:O	1:A:249:LYS:HG2	1.95	0.65
1:A:447:TRP:CE3	1:A:451:ILE:HG21	2.31	0.65
1:A:66:VAL:HG12	1:A:67:GLY:N	2.11	0.65
1:A:86:ASP:O	1:A:87:GLY:C	2.35	0.64
1:A:249:LYS:CD	2:A:589:MPD:C5	2.75	0.64
1:A:97:GLU:HB3	1:A:105:ARG:O	1.96	0.64
1:A:461:LYS:NZ	2:A:591:MPD:HM1	2.10	0.64
1:A:254:GLN:NE2	1:A:325:ASP:H	1.95	0.64
1:A:102:PHE:O	1:A:103:ILE:HD12	1.95	0.64
1:A:87:GLY:N	1:A:96:ARG:HH22	1.96	0.64
1:A:101:ASP:OD1	1:A:102:PHE:N	2.31	0.64
1:A:261:CYS:CB	1:A:332:SER:HB2	2.25	0.64
1:A:97:GLU:C	1:A:97:GLU:OE1	2.35	0.64
1:A:487:ARG:HD2	4:A:670:HOH:O	1.98	0.63
1:A:233:PRO:HB3	1:A:392:VAL:CG2	2.22	0.63
1:A:585:LYS:O	1:A:586:TYR:HB2	1.98	0.63
1:A:99:THR:CG2	1:A:100:GLY:N	2.49	0.63
1:A:401:GLU:HG2	1:A:402:PRO:CD	2.26	0.63
1:A:458:MET:SD	1:A:525:ARG:HG2	2.39	0.63
1:A:528:LEU:HB3	2:A:591:MPD:HM2	1.81	0.62
1:A:80:LEU:HD21	1:A:89:GLN:OE1	1.99	0.62
1:A:102:PHE:O	1:A:103:ILE:CD1	2.47	0.62
1:A:102:PHE:O	1:A:103:ILE:HG13	1.99	0.62
1:A:96:ARG:CG	1:A:301:PRO:CD	2.68	0.61
1:A:200:PRO:HG2	1:A:377:LEU:HD11	1.81	0.61
1:A:249:LYS:HZ2	1:A:281:PRO:HG3	1.64	0.61
1:A:96:ARG:NH2	1:A:304:ALA:HB1	2.15	0.61
1:A:258:TYR:CE2	1:A:260:GLY:HA2	2.34	0.60
1:A:329:MET:HA	4:A:838:HOH:O	2.01	0.60
1:A:394:PRO:HD2	4:A:766:HOH:O	2.01	0.60
1:A:494:VAL:HG11	1:A:531:TYR:HB2	1.83	0.60
1:A:86:ASP:HB2	1:A:96:ARG:CZ	2.31	0.60
1:A:109:ALA:HB1	1:A:110:PRO:CD	2.31	0.60
1:A:448:HIS:CA	1:A:452:ASP:O	2.47	0.59
1:A:86:ASP:O	1:A:88:ILE:N	2.35	0.59
1:A:67:GLY:CA	4:A:717:HOH:O	2.50	0.59
1:A:540:ARG:NH2	1:A:588:ALA:OXT	2.36	0.59
1:A:232:ILE:HG12	1:A:232:ILE:O	1.99	0.58
1:A:189:MET:SD	4:A:756:HOH:O	2.57	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:233:PRO:CB	1:A:392:VAL:HG23	2.24	0.58
1:A:394:PRO:HB2	1:A:398:ASP:HB3	1.85	0.58
1:A:502:TYR:OH	1:A:520:GLN:HB3	2.02	0.58
1:A:532:ASP:OD2	2:A:591:MPD:HM3	2.04	0.58
1:A:102:PHE:O	1:A:103:ILE:CG1	2.51	0.58
1:A:90:ARG:NH2	1:A:95:ILE:O	2.37	0.58
1:A:290:ARG:CZ	4:A:804:HOH:O	2.52	0.58
1:A:199:ARG:NH1	1:A:318:TYR:HA	2.18	0.58
1:A:245:HIS:CD2	2:A:589:MPD:C3	2.83	0.57
1:A:197:VAL:HA	4:A:652:HOH:O	2.05	0.57
1:A:90:ARG:HH22	1:A:96:ARG:CA	2.14	0.57
1:A:103:ILE:CG2	1:A:104:ALA:H	2.17	0.57
1:A:554:GLU:OE2	1:A:558:ARG:NH2	2.32	0.57
1:A:587:GLY:O	1:A:588:ALA:HB3	2.04	0.57
1:A:364:LEU:HD21	1:A:407:THR:OG1	2.05	0.57
1:A:118:TRP:HH2	4:A:818:HOH:O	1.86	0.56
1:A:331:ASP:HA	1:A:391:ALA:HB3	1.88	0.56
1:A:203:GLU:OE2	4:A:696:HOH:O	2.17	0.55
1:A:250:TRP:CH2	1:A:281:PRO:HB2	2.40	0.55
1:A:245:HIS:CD2	2:A:589:MPD:C5	2.86	0.55
1:A:451:ILE:HG22	1:A:521:VAL:HG11	1.89	0.55
1:A:65:VAL:HG12	1:A:66:VAL:O	2.07	0.55
1:A:245:HIS:HE1	1:A:278:LEU:HD13	1.71	0.55
1:A:75:GLU:HB2	1:A:112:LEU:HD13	1.89	0.55
1:A:259:ILE:C	1:A:261:CYS:SG	2.85	0.55
1:A:448:HIS:CE1	1:A:456:LYS:H	2.22	0.55
1:A:213:THR:HB	1:A:219:ASP:OD1	2.06	0.54
1:A:110:PRO:HG2	4:A:819:HOH:O	2.05	0.54
1:A:577:GLU:HB3	4:A:793:HOH:O	2.08	0.54
1:A:226:LYS:NZ	1:A:252:ASP:OD2	2.21	0.54
1:A:108:THR:CG2	1:A:109:ALA:H	2.16	0.54
1:A:153:PRO:HG2	1:A:193:TRP:HZ3	1.72	0.53
1:A:421:LEU:HD13	2:A:590:MPD:O4	2.09	0.53
1:A:458:MET:HE1	1:A:525:ARG:HA	1.91	0.53
1:A:249:LYS:NZ	1:A:249:LYS:HB2	2.24	0.53
1:A:261:CYS:HB2	1:A:332:SER:OG	2.09	0.53
1:A:417:LEU:HA	1:A:429:ALA:O	2.08	0.52
1:A:474:GLU:OE2	1:A:474:GLU:HA	2.08	0.52
1:A:170:ILE:HG12	4:A:657:HOH:O	2.09	0.52
1:A:258:TYR:HE2	1:A:260:GLY:CA	2.05	0.52
1:A:96:ARG:O	1:A:301:PRO:HG2	2.10	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:249:LYS:HD3	2:A:589:MPD:C5	2.40	0.52
1:A:458:MET:HE1	1:A:525:ARG:HG2	1.89	0.51
1:A:232:ILE:O	1:A:232:ILE:CG1	2.59	0.51
1:A:86:ASP:OD2	1:A:90:ARG:O	2.29	0.51
1:A:90:ARG:HD3	1:A:108:THR:HG23	1.92	0.51
1:A:108:THR:CG2	1:A:109:ALA:N	2.74	0.51
1:A:501:ASP:O	1:A:560:LYS:HE2	2.11	0.51
1:A:452:ASP:OD2	1:A:525:ARG:NH1	2.44	0.51
1:A:86:ASP:OD1	1:A:96:ARG:NH2	2.44	0.51
1:A:96:ARG:CB	1:A:301:PRO:CD	2.83	0.51
1:A:447:TRP:HZ3	1:A:451:ILE:HG21	1.75	0.50
1:A:585:LYS:O	1:A:586:TYR:CB	2.59	0.50
1:A:468:LYS:HE3	1:A:472:LEU:HG	1.93	0.50
1:A:66:VAL:CG1	1:A:67:GLY:N	2.74	0.50
1:A:199:ARG:HG3	1:A:317:GLU:OE1	2.11	0.50
1:A:90:ARG:HH11	1:A:90:ARG:HG3	1.76	0.50
1:A:373:ARG:HG3	1:A:385:SER:HB3	1.93	0.50
1:A:236:PHE:CE1	1:A:240:LYS:HE2	2.46	0.50
1:A:79:ARG:CG	4:A:818:HOH:O	1.85	0.49
1:A:249:LYS:CE	2:A:589:MPD:H32	2.42	0.49
1:A:75:GLU:HB2	1:A:112:LEU:CD1	2.42	0.49
1:A:97:GLU:OE1	1:A:97:GLU:O	2.30	0.49
1:A:209:VAL:HG22	1:A:446:TRP:CD1	2.47	0.49
1:A:451:ILE:HD13	1:A:451:ILE:N	2.27	0.49
1:A:79:ARG:CB	1:A:118:TRP:HZ2	2.04	0.49
1:A:249:LYS:HD3	2:A:589:MPD:H52	1.93	0.49
1:A:452:ASP:OD1	1:A:453:PRO:CD	2.60	0.49
1:A:88:ILE:N	1:A:88:ILE:CD1	2.75	0.49
1:A:582:LEU:O	1:A:585:LYS:O	2.31	0.49
1:A:97:GLU:OE1	1:A:98:LYS:O	2.31	0.48
1:A:536:GLU:O	1:A:539:ASN:ND2	2.47	0.48
1:A:74:VAL:HG23	1:A:195:VAL:HG11	1.93	0.48
1:A:147:HIS:HE1	1:A:318:TYR:OH	1.96	0.48
1:A:290:ARG:NH2	4:A:804:HOH:O	2.45	0.48
1:A:258:TYR:N	4:A:851:HOH:O	2.36	0.48
1:A:288:MET:CE	1:A:293:LEU:CD1	2.92	0.48
1:A:207:PRO:O	1:A:442:ALA:HB1	2.13	0.48
1:A:102:PHE:CB	1:A:103:ILE:HD12	2.39	0.48
1:A:241:THR:OG1	1:A:264:ARG:NH2	2.47	0.48
1:A:114:ARG:HD3	1:A:170:ILE:HD11	1.96	0.48
1:A:176:LYS:HE2	4:A:857:HOH:O	2.14	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:102:PHE:HB2	1:A:103:ILE:CD1	2.39	0.47
1:A:458:MET:HE3	1:A:525:ARG:HG2	1.95	0.47
1:A:67:GLY:HA2	4:A:717:HOH:O	2.13	0.47
1:A:209:VAL:HG13	1:A:210:PRO:HD2	1.97	0.47
1:A:66:VAL:HG12	1:A:67:GLY:H	1.78	0.47
1:A:258:TYR:HB3	1:A:293:LEU:HD23	1.96	0.47
1:A:499:ARG:HA	1:A:503:LEU:HB2	1.95	0.47
1:A:90:ARG:HD2	1:A:94:VAL:HB	1.96	0.47
1:A:364:LEU:CD2	1:A:407:THR:OG1	2.62	0.47
1:A:578:GLN:HA	1:A:581:GLU:HG2	1.96	0.47
1:A:393:SER:HB3	4:A:813:HOH:O	2.13	0.47
1:A:538:ILE:C	1:A:540:ARG:H	2.17	0.47
1:A:380:ASP:OD2	1:A:382:ARG:HD3	2.15	0.47
1:A:103:ILE:HG23	1:A:104:ALA:H	1.80	0.47
1:A:261:CYS:HB3	1:A:296:ASN:N	2.25	0.47
2:A:589:MPD:O4	2:A:589:MPD:CM	2.62	0.47
1:A:452:ASP:OD1	1:A:453:PRO:HD2	2.15	0.47
1:A:249:LYS:HE2	2:A:589:MPD:H32	1.96	0.46
1:A:545:GLU:O	1:A:549:LYS:HG2	2.15	0.46
1:A:216:ARG:HD2	1:A:502:TYR:CZ	2.50	0.46
1:A:202:LYS:NZ	4:A:835:HOH:O	2.20	0.46
1:A:414:PHE:CE2	1:A:416:ALA:HB2	2.51	0.46
1:A:449:LYS:HE2	4:A:840:HOH:O	2.16	0.46
1:A:399:PHE:O	1:A:399:PHE:CD2	2.69	0.46
1:A:257:ILE:HB	1:A:328:LEU:HD12	1.98	0.46
1:A:79:ARG:CB	1:A:118:TRP:CZ2	2.77	0.45
1:A:302:VAL:HG21	1:A:338:GLU:OE2	2.16	0.45
1:A:458:MET:HE2	2:A:591:MPD:H31	1.98	0.45
1:A:141:GLU:OE1	1:A:147:HIS:HD2	1.98	0.45
1:A:88:ILE:H	1:A:96:ARG:NH2	2.14	0.45
1:A:288:MET:HE2	1:A:293:LEU:CD1	2.47	0.45
1:A:97:GLU:O	1:A:303:ALA:HB3	2.16	0.45
1:A:233:PRO:HA	1:A:392:VAL:O	2.17	0.45
1:A:97:GLU:O	1:A:303:ALA:CB	2.65	0.45
1:A:501:ASP:OD2	1:A:556:ILE:HG22	2.17	0.45
1:A:262:GLY:O	1:A:335:ARG:NE	2.48	0.45
1:A:83:SER:OG	1:A:85:TYR:HE2	1.87	0.44
1:A:102:PHE:C	1:A:103:ILE:CD1	2.78	0.44
1:A:432:TRP:CH2	2:A:590:MPD:HM3	2.52	0.44
1:A:233:PRO:CD	1:A:414:PHE:HE1	2.01	0.44
1:A:336:TRP:CH2	1:A:364:LEU:HD13	2.52	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:245:HIS:NE2	1:A:249:LYS:HD3	2.27	0.44
1:A:240:LYS:HE3	1:A:240:LYS:HB3	1.79	0.44
1:A:71:SER:C	4:A:808:HOH:O	2.55	0.44
1:A:79:ARG:HA	4:A:802:HOH:O	2.17	0.44
1:A:583:PHE:HB3	1:A:588:ALA:O	2.18	0.44
2:A:590:MPD:O4	2:A:590:MPD:CM	2.65	0.44
1:A:66:VAL:CG1	1:A:67:GLY:H	2.31	0.44
1:A:331:ASP:OD1	1:A:331:ASP:O	2.36	0.44
1:A:249:LYS:HB2	1:A:249:LYS:HZ3	1.83	0.43
1:A:587:GLY:O	1:A:588:ALA:CB	2.67	0.43
1:A:60:ARG:HG3	1:A:108:THR:HG21	2.00	0.43
1:A:94:VAL:O	1:A:96:ARG:HD2	2.18	0.43
1:A:77:GLY:H	1:A:80:LEU:HD12	1.82	0.43
1:A:198:LYS:NZ	4:A:732:HOH:O	2.48	0.43
1:A:420:ASP:OD1	1:A:423:ARG:NH1	2.52	0.43
1:A:261:CYS:HB2	1:A:296:ASN:HB2	2.00	0.43
1:A:366:GLU:HA	1:A:369:GLU:HB2	2.01	0.43
1:A:74:VAL:HG23	1:A:195:VAL:HG12	1.98	0.43
1:A:72:LEU:HD12	1:A:196:ARG:HH12	1.84	0.42
1:A:76:LEU:HA	1:A:80:LEU:HD11	2.01	0.42
1:A:536:GLU:O	1:A:540:ARG:CG	2.66	0.42
1:A:130:VAL:HG11	1:A:159:ILE:HD11	2.01	0.42
1:A:171:GLU:O	1:A:188:LYS:HE2	2.19	0.42
1:A:212:ILE:HD13	1:A:212:ILE:HA	1.89	0.42
1:A:244:GLN:HE22	1:A:391:ALA:HB2	1.84	0.42
1:A:338:GLU:CG	1:A:339:ALA:N	2.82	0.42
1:A:92:LEU:HD23	1:A:92:LEU:O	2.18	0.42
1:A:204:LYS:HD2	1:A:372:GLY:HA3	2.02	0.42
3:A:592:TRS:H31	4:A:856:HOH:O	2.18	0.42
1:A:161:GLU:OE1	1:A:176:LYS:HE3	2.20	0.42
1:A:118:TRP:O	1:A:167:ASP:HA	2.20	0.42
1:A:216:ARG:N	1:A:505:GLN:HE22	2.00	0.42
1:A:548:ALA:O	1:A:553:ARG:NH1	2.51	0.42
1:A:78:PRO:O	1:A:79:ARG:CB	2.67	0.42
1:A:291:THR:CG2	4:A:851:HOH:O	2.68	0.42
1:A:301:PRO:O	1:A:302:VAL:C	2.57	0.41
1:A:86:ASP:C	1:A:88:ILE:N	2.71	0.41
1:A:315:ILE:HD13	1:A:315:ILE:HG21	1.86	0.41
1:A:525:ARG:HB3	4:A:841:HOH:O	2.20	0.41
1:A:584:LYS:HB2	1:A:584:LYS:HE2	1.86	0.41
1:A:70:ALA:O	1:A:71:SER:CB	2.68	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:287:LEU:HD12	1:A:287:LEU:HA	1.88	0.41
1:A:577:GLU:HG3	4:A:791:HOH:O	2.20	0.41
1:A:426:HIS:HD2	4:A:677:HOH:O	2.02	0.41
1:A:271:VAL:HA	1:A:275:PHE:HD1	1.86	0.41
1:A:447:TRP:CE3	1:A:451:ILE:HG13	2.56	0.41
1:A:546:GLU:CG	1:A:582:LEU:HD11	2.39	0.41
1:A:73:SER:HB2	1:A:193:TRP:O	2.20	0.41
1:A:88:ILE:H	1:A:96:ARG:HH22	1.69	0.41
1:A:261:CYS:CB	1:A:332:SER:CB	2.93	0.41
1:A:288:MET:CE	1:A:293:LEU:HD11	2.50	0.41
1:A:114:ARG:NH1	1:A:190:TYR:CZ	2.78	0.41
1:A:108:THR:HG22	1:A:109:ALA:O	2.21	0.40
1:A:69:GLY:CA	1:A:103:ILE:HG13	2.47	0.40
1:A:87:GLY:HA2	1:A:308:SER:CB	2.45	0.40
1:A:486:GLU:HB3	1:A:538:ILE:HD13	2.03	0.40
1:A:249:LYS:CD	2:A:589:MPD:H52	2.49	0.40
1:A:97:GLU:C	1:A:97:GLU:CD	2.80	0.40
1:A:114:ARG:O	1:A:169:THR:HB	2.22	0.40
1:A:193:TRP:CD2	1:A:194:PRO:HD2	2.57	0.40
1:A:394:PRO:HA	1:A:395:PRO:HD3	1.85	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	509/588 (87%)	465 (91%)	35 (7%)	9 (2%)	8 10

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	71	SER

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Mol	Chain	Res	Type
1	A	237	GLY
1	A	262	GLY
1	A	233	PRO
1	A	397	GLY
1	A	401	GLU
1	A	586	TYR
1	A	403	VAL
1	A	88	ILE

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	434/492 (88%)	390 (90%)	44 (10%)	7 11

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	60	ARG
1	A	68	THR
1	A	72	LEU
1	A	83	SER
1	A	88	ILE
1	A	92	LEU
1	A	94	VAL
1	A	95	ILE
1	A	96	ARG
1	A	97	GLU
1	A	98	LYS
1	A	105	ARG
1	A	129	LYS
1	A	181	SER
1	A	204	LYS
1	A	213	THR
1	A	232	ILE
1	A	236	PHE

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Mol	Chain	Res	Type
1	A	240	LYS
1	A	242	VAL
1	A	249	LYS
1	A	261	CYS
1	A	288	MET
1	A	290	ARG
1	A	298	SER
1	A	299	ASN
1	A	305	ARG
1	A	326	VAL
1	A	363	LYS
1	A	366	GLU
1	A	369	GLU
1	A	404	VAL
1	A	406	ASN
1	A	408	LEU
1	A	417	LEU
1	A	421	LEU
1	A	433	LEU
1	A	450	ASN
1	A	459	ARG
1	A	474	GLU
1	A	493	LEU
1	A	539	ASN
1	A	540	ARG
1	A	580	GLU

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	89	GLN
1	A	147	HIS
1	A	245	HIS
1	A	254	GLN
1	A	266	ASN
1	A	299	ASN
1	A	426	HIS
1	A	448	HIS
1	A	450	ASN
1	A	504	GLN
1	A	505	GLN
1	A	539	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

4 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	MPD	A	591	-	7,7,7	1.24	2 (28%)	9,10,10	1.12	1 (11%)
2	MPD	A	590	-	7,7,7	1.23	2 (28%)	9,10,10	0.30	0
3	TRS	A	592	-	7,7,7	0.27	0	9,9,9	0.53	0
2	MPD	A	589	-	7,7,7	1.24	2 (28%)	9,10,10	0.22	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	MPD	A	591	-	-	2/5/5/5	-
2	MPD	A	590	-	-	2/5/5/5	-
3	TRS	A	592	-	-	2/9/9/9	-
2	MPD	A	589	-	-	0/5/5/5	-

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	591	MPD	O4-C4	2.34	1.53	1.43
2	A	590	MPD	O4-C4	2.34	1.53	1.43
2	A	589	MPD	O4-C4	2.29	1.53	1.43
2	A	589	MPD	C5-C4	-2.28	1.41	1.51
2	A	590	MPD	C5-C4	-2.21	1.41	1.51
2	A	591	MPD	C5-C4	-2.16	1.41	1.51

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	591	MPD	CM-C2-C1	-2.35	105.67	110.57

There are no chirality outliers.

All (6) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	590	MPD	CM-C2-C3-C4
2	A	591	MPD	C1-C2-C3-C4
3	A	592	TRS	C2-C-C3-O3
2	A	590	MPD	O2-C2-C3-C4
2	A	591	MPD	O2-C2-C3-C4
3	A	592	TRS	N-C-C3-O3

There are no ring outliers.

4 monomers are involved in 28 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	591	MPD	6	0
2	A	590	MPD	3	0
3	A	592	TRS	1	0
2	A	589	MPD	18	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	513/588 (87%)	0.67	77 (15%) 2 1	26, 52, 104, 123	0

All (77) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	395	PRO	11.0
1	A	96	ARG	10.6
1	A	62	GLY	9.7
1	A	396	GLY	8.7
1	A	60	ARG	7.5
1	A	100	GLY	7.4
1	A	61	PRO	7.2
1	A	408	LEU	6.8
1	A	64	PRO	6.4
1	A	397	GLY	6.3
1	A	63	GLU	6.1
1	A	588	ALA	5.8
1	A	95	ILE	5.7
1	A	108	THR	5.5
1	A	405	GLN	5.4
1	A	584	LYS	5.2
1	A	399	PHE	5.2
1	A	336	TRP	5.1
1	A	66	VAL	5.1
1	A	404	VAL	5.0
1	A	99	THR	5.0
1	A	107	VAL	5.0
1	A	101	ASP	4.9
1	A	273	GLU	4.5
1	A	238	ALA	4.3
1	A	402	PRO	4.2
1	A	502	TYR	4.2

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Mol	Chain	Res	Type	RSRZ
1	A	65	VAL	4.2
1	A	299	ASN	4.1
1	A	294	ILE	3.9
1	A	401	GLU	3.9
1	A	85	TYR	3.9
1	A	269	THR	3.9
1	A	407	THR	3.6
1	A	406	ASN	3.6
1	A	361	ALA	3.4
1	A	266	ASN	3.4
1	A	359	TYR	3.3
1	A	400	SER	3.3
1	A	102	PHE	3.2
1	A	264	ARG	3.2
1	A	237	GLY	3.1
1	A	389	ILE	3.0
1	A	360	LEU	3.0
1	A	89	GLN	3.0
1	A	103	ILE	2.9
1	A	257	ILE	2.9
1	A	311	THR	2.9
1	A	328	LEU	2.9
1	A	358	ALA	2.9
1	A	268	MET	2.8
1	A	403	VAL	2.8
1	A	394	PRO	2.8
1	A	80	LEU	2.7
1	A	270	ASP	2.7
1	A	208	GLU	2.7
1	A	398	ASP	2.7
1	A	330	ALA	2.6
1	A	293	LEU	2.6
1	A	106	GLY	2.6
1	A	235	PRO	2.6
1	A	451	ILE	2.6
1	A	259	ILE	2.6
1	A	452	ASP	2.4
1	A	213	THR	2.4
1	A	272	LEU	2.4
1	A	298	SER	2.3
1	A	309	ILE	2.3
1	A	327	ALA	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	587	GLY	2.3
1	A	94	VAL	2.3
1	A	70	ALA	2.2
1	A	76	LEU	2.2
1	A	84	ILE	2.1
1	A	581	GLU	2.1
1	A	329	MET	2.1
1	A	88	ILE	2.0

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

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

6.3 Carbohydrates (i)

There are no monosaccharides 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(Å^2)	Q<0.9
2	MPD	A	589	8/8	0.70	0.37	61,63,64,67	0
2	MPD	A	591	8/8	0.73	0.28	74,76,77,78	0
3	TRS	A	592	8/8	0.76	0.20	93,95,95,95	0
2	MPD	A	590	8/8	0.83	0.29	76,79,80,80	0

6.5 Other polymers (i)

There are no such residues in this entry.