



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:04 AM GMT

PDB ID : 3KRO
Title : Mint heterotetrameric geranyl pyrophosphate synthase in complex with magnesium, IPP, and DMASPP (II)
Authors : Chang, T.-H.; Hsieh, F.-L.; Ko, T.-P.; Wang, A.H.-J.
Deposited on : 2009-11-19
Resolution : 1.95 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

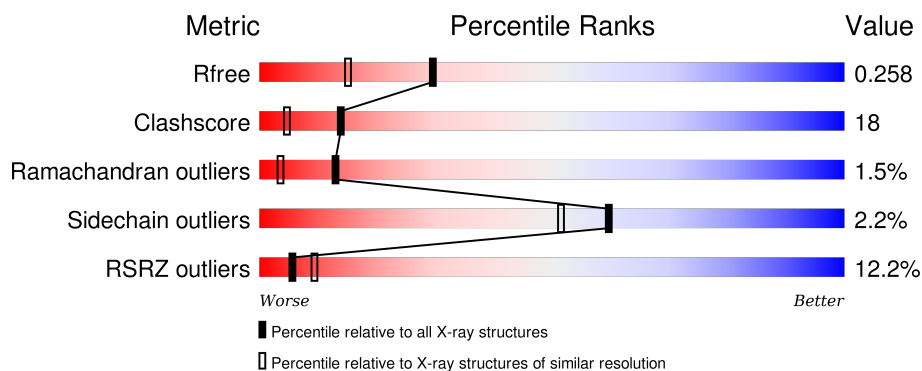
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.95 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	1833 (1.96-1.96)
Clashscore	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)
RSRZ outliers	91569	1835 (1.96-1.96)

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	295	<div> <div>7%</div> <div>74%</div> <div>20%</div> <div>• 5%</div> </div>
1	D	295	<div> <div>8%</div> <div>73%</div> <div>26%</div> <div>•</div> </div>
2	B	274	<div> <div>18%</div> <div>58%</div> <div>33%</div> <div>• • 5%</div> </div>
2	C	274	<div> <div>14%</div> <div>65%</div> <div>26%</div> <div>• 5%</div> </div>

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 crite-

ria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	EDO	C	2005	-	-	-	X

2 Entry composition

There are 8 unique types of molecules in this entry. The entry contains 9441 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 Geranyl diphosphate synthase large subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	279	Total	C	N	O	S	0	0	0
			2120	1338	372	392	18			
1	D	295	Total	C	N	O	S	0	0	0
			2232	1406	391	417	18			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MET	-	EXPRESSION TAG	UNP Q9SBR3
D	1	MET	-	EXPRESSION TAG	UNP Q9SBR3

- Molecule 2 is a protein called Geranyl diphosphate synthase small subunit.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	259	Total	C	N	O	S	0	0	0
			1954	1233	339	368	14			
2	C	259	Total	C	N	O	S	0	0	0
			1954	1233	339	368	14			

There are 18 discrepancies between the modelled and reference sequences:

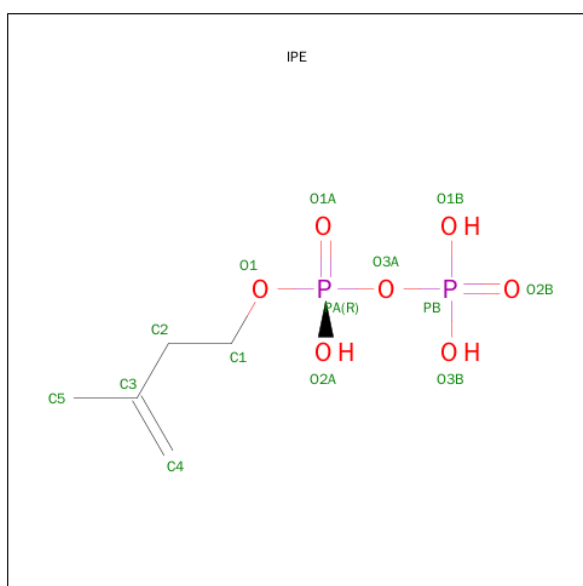
Chain	Residue	Modelled	Actual	Comment	Reference
B	1	MET	-	EXPRESSION TAG	UNP Q9SBR4
B	267	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	268	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	269	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	270	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	271	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	272	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	273	HIS	-	EXPRESSION TAG	UNP Q9SBR4
B	274	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	1	MET	-	EXPRESSION TAG	UNP Q9SBR4

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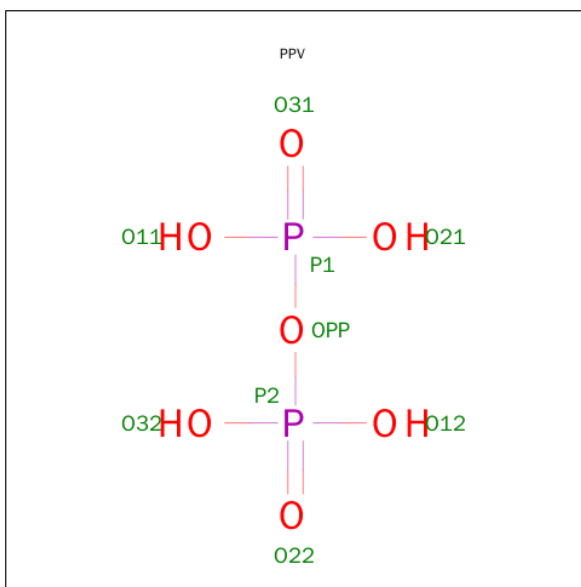
Chain	Residue	Modelled	Actual	Comment	Reference
C	267	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	268	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	269	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	270	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	271	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	272	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	273	HIS	-	EXPRESSION TAG	UNP Q9SBR4
C	274	HIS	-	EXPRESSION TAG	UNP Q9SBR4

- Molecule 3 is 3-METHYLBUT-3-ENYL TRIHYDROGEN DIPHOSPHATE (three-letter code: IPE) (formula: $C_5H_{12}O_7P_2$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	O	P	0	0
			14	5	7	2		
3	D	1	Total	C	O	P	0	0
			14	5	7	2		

- Molecule 4 is PYROPHOSPHATE (three-letter code: PPV) (formula: $H_4O_7P_2$).

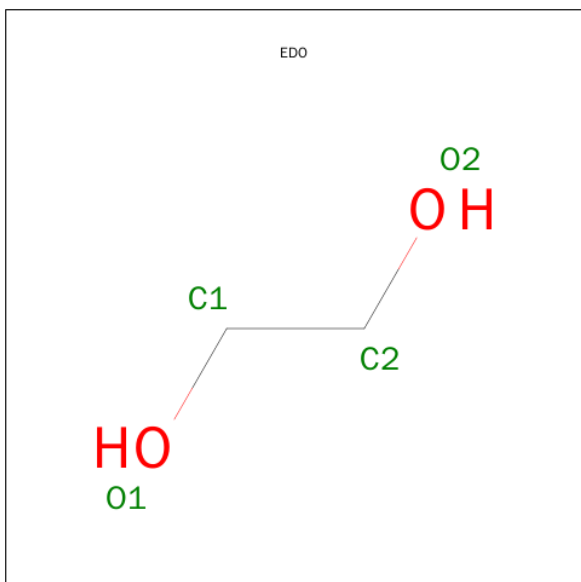


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O P 9 7 2	0	0

- Molecule 5 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

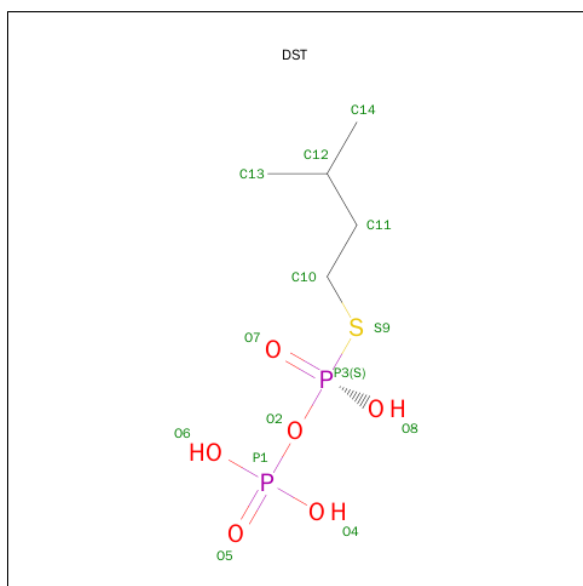
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	2	Total Mg 2 2	0	0
5	D	2	Total Mg 2 2	0	0

- Molecule 6 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	C	1	Total	C	O	0	0
			4	2	2		

- Molecule 7 is DIMETHYLALLYL S-THIOLODIPHOSPHATE (three-letter code: DST) (formula: $C_5H_{14}O_6P_2S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
7	D	1	Total	C	O	P	S	0	0
			14	5	6	2	1		

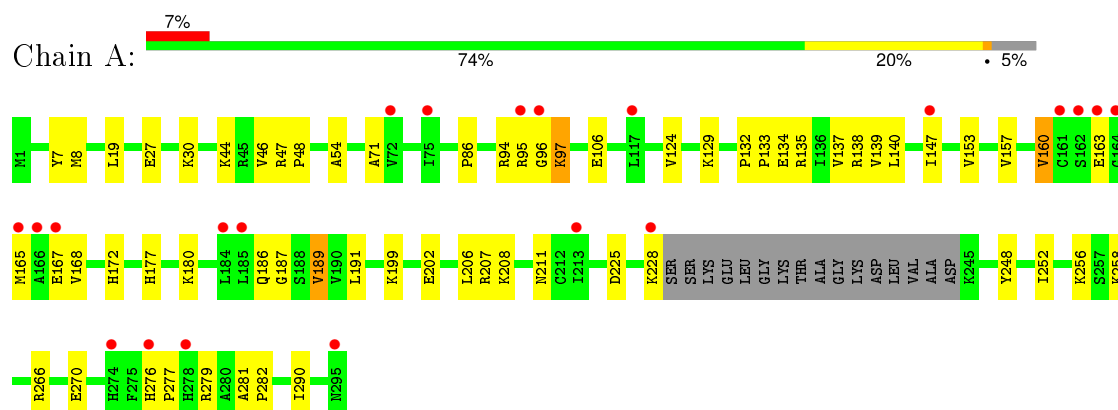
- Molecule 8 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
8	A	268	Total	O	0	0
			268	268		
8	B	243	Total	O	0	0
			243	243		
8	C	294	Total	O	0	0
			294	294		
8	D	317	Total	O	0	0
			317	317		

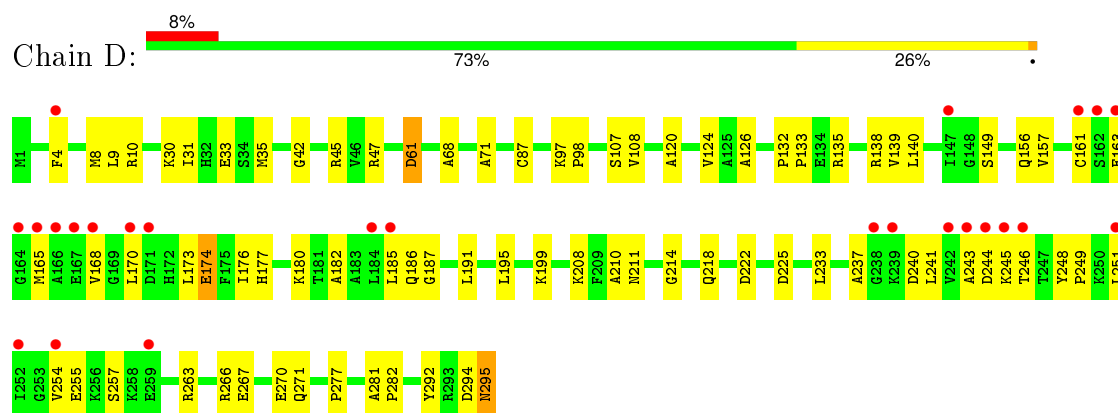
3 Residue-property plots

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.

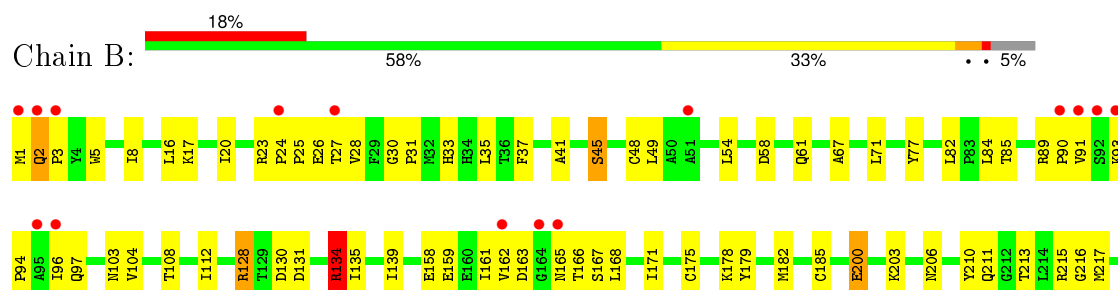
- Molecule 1: Geranyl diphosphate synthase large subunit

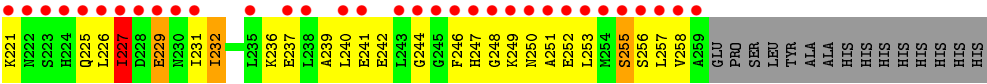


- Molecule 1: Geranyl diphosphate synthase large subunit

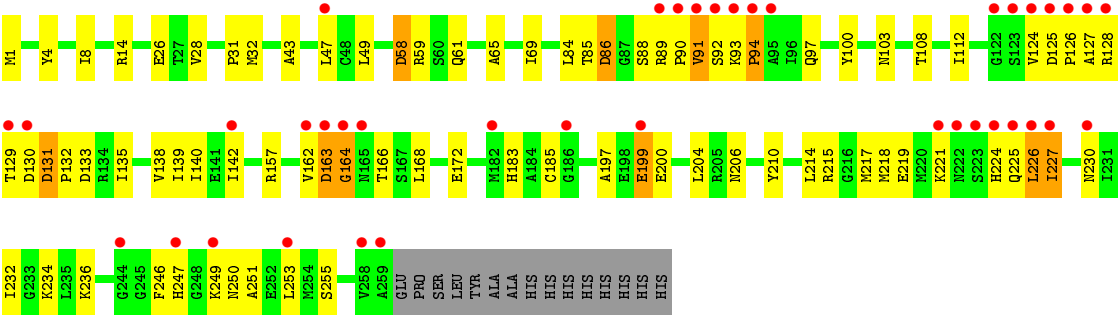


- Molecule 2: Geranyl diphosphate synthase small subunit





● Molecule 2: Geranyl diphosphate synthase small subunit



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	53.44Å 108.70Å 182.28Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 – 1.95 29.02 – 1.94	Depositor EDS
% Data completeness (in resolution range)	90.1 (30.00-1.95) 89.6 (29.02-1.94)	Depositor EDS
R_{merge}	0.05	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.34 (at 1.95Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.205 , 0.256 0.205 , 0.258	Depositor DCC
R_{free} test set	3588 reflections (5.08%)	DCC
Wilson B-factor (Å ²)	30.2	Xtriage
Anisotropy	0.820	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 57.1	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	1 of 70905 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	9441	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.96% 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, 0.375 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: PPV, MG, EDO, DST, IPE

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	2/2155 (0.1%)	0.85	0/2903
1	D	0.93	0/2268	0.87	3/3055 (0.1%)
2	B	0.83	2/1993 (0.1%)	0.82	1/2695 (0.0%)
2	C	0.84	1/1993 (0.1%)	0.83	1/2695 (0.0%)
All	All	0.89	5/8409 (0.1%)	0.84	5/11348 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	175	CYS	CB-SG	5.74	1.92	1.82
1	A	27	GLU	CG-CD	5.23	1.59	1.51
1	A	46	VAL	CA-CB	5.20	1.65	1.54
2	B	200	GLU	CG-CD	5.15	1.59	1.51
2	C	185	CYS	CB-SG	-5.05	1.73	1.81

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	47	ARG	NE-CZ-NH2	-6.67	116.96	120.30
2	C	157	ARG	NE-CZ-NH1	-6.08	117.26	120.30
1	D	45	ARG	NE-CZ-NH2	-5.75	117.43	120.30
1	D	61	ASP	CB-CG-OD2	-5.44	113.41	118.30
2	B	134	ARG	NE-CZ-NH1	5.32	122.96	120.30

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	2120	0	2146	51	0
1	D	2232	0	2264	79	0
2	B	1954	0	1948	103	0
2	C	1954	0	1948	83	0
3	A	14	0	9	1	0
3	D	14	0	9	0	0
4	A	9	0	0	0	0
5	A	2	0	0	0	0
5	D	2	0	0	0	0
6	C	4	0	6	0	0
7	D	14	0	10	3	0
8	A	268	0	0	14	1
8	B	243	0	0	13	1
8	C	294	0	0	15	1
8	D	317	0	0	15	1
All	All	9441	0	8340	307	2

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 18.

All (307) 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)
2:B:2:GLN:HB2	2:B:3:PRO:HD3	1.21	1.20
2:B:229:GLU:HG3	8:B:289:HOH:O	1.39	1.17
2:C:250:ASN:HB3	8:C:1039:HOH:O	1.57	1.03
2:C:230:ASN:O	2:C:234:LYS:HG3	1.58	1.02
2:C:28:VAL:HG22	1:D:157:VAL:CG2	1.96	0.94
2:C:61:GLN:HE21	2:C:128:ARG:HH22	0.94	0.93
2:B:221:LYS:HG3	2:B:227:ILE:HG12	1.47	0.93
1:D:4:PHE:HB2	1:D:8:MET:HE3	1.51	0.92
1:D:208:LYS:HE2	8:D:1010:HOH:O	1.70	0.90
1:D:10:ARG:HD3	8:D:583:HOH:O	1.72	0.88
2:B:58:ASP:H	2:B:61:GLN:HE21	1.21	0.88
2:C:61:GLN:NE2	2:C:128:ARG:HH22	1.72	0.87

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:28:VAL:HG22	1:D:157:VAL:HG21	1.56	0.87
2:B:25:PRO:HB2	2:B:27:THR:HG22	1.57	0.86
1:D:4:PHE:HB2	1:D:8:MET:CE	2.06	0.85
2:B:26:GLU:HB3	8:B:435:HOH:O	1.75	0.85
2:B:2:GLN:HB2	2:B:3:PRO:CD	2.04	0.84
1:A:97:LYS:HD2	1:A:97:LYS:N	1.91	0.84
2:B:240:LEU:HD21	2:B:258:VAL:HG21	1.59	0.84
2:B:227:ILE:HG23	8:B:290:HOH:O	1.77	0.83
2:B:5:TRP:HZ2	2:B:253:LEU:HG	1.42	0.83
2:B:249:LYS:HD3	2:B:253:LEU:HB2	1.59	0.83
2:B:94:PRO:HB2	8:B:968:HOH:O	1.79	0.82
2:B:5:TRP:CZ2	2:B:253:LEU:HG	2.13	0.82
2:C:221:LYS:HB2	2:C:227:ILE:HD12	1.61	0.82
2:B:23:ARG:HB2	2:B:24:PRO:HD2	1.62	0.81
2:C:221:LYS:HD3	2:C:227:ILE:HB	1.62	0.80
2:B:162:VAL:H	2:B:166:THR:CG2	1.95	0.80
2:C:247:HIS:HB3	8:C:1129:HOH:O	1.80	0.80
1:A:163:GLU:HG2	8:A:702:HOH:O	1.80	0.79
2:C:61:GLN:HE21	2:C:128:ARG:NH2	1.78	0.79
2:B:229:GLU:OE1	2:B:232:ILE:HD11	1.81	0.78
2:B:77:TYR:OH	2:B:96:ILE:HD13	1.83	0.78
2:B:128:ARG:CG	2:B:128:ARG:HH11	1.96	0.78
2:B:253:LEU:HD12	2:B:256:SER:OG	1.84	0.77
2:B:71:LEU:HD22	2:B:112:ILE:HG23	1.67	0.77
1:D:173:LEU:HD13	1:D:246:THR:HG22	1.64	0.77
2:C:8:ILE:HG22	2:C:49:LEU:HD12	1.66	0.76
2:B:249:LYS:O	2:B:253:LEU:N	2.18	0.76
2:B:206:ASN:CB	2:B:242:GLU:HG2	2.15	0.76
2:C:214:LEU:HD11	2:C:232:ILE:HG23	1.69	0.75
1:A:168:VAL:HG13	1:A:172:HIS:HB3	1.68	0.75
2:B:128:ARG:HG2	2:B:128:ARG:HH11	1.52	0.75
1:D:170:LEU:HD23	1:D:251:LEU:HD13	1.69	0.74
2:B:206:ASN:HB2	2:B:242:GLU:HG2	1.70	0.74
1:A:256:LYS:HE2	8:A:547:HOH:O	1.87	0.74
2:C:227:ILE:CG1	8:C:1126:HOH:O	2.35	0.74
2:C:250:ASN:HA	2:C:253:LEU:HD12	1.70	0.73
2:C:93:LYS:N	2:C:94:PRO:HD2	2.04	0.73
1:A:134:GLU:HG3	8:A:930:HOH:O	1.88	0.72
2:B:97:GLN:OE1	8:B:851:HOH:O	2.09	0.71
2:B:162:VAL:H	2:B:166:THR:HG21	1.57	0.70
1:D:132:PRO:HG2	1:D:135:ARG:HB2	1.74	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:135:ARG:HG3	1:D:135:ARG:HH11	1.55	0.70
2:B:250:ASN:HA	2:B:253:LEU:HB3	1.74	0.70
1:D:163:GLU:O	1:D:244:ASP:HB3	1.92	0.69
1:D:68:ALA:HA	1:D:124:VAL:HG22	1.73	0.69
1:D:180:LYS:HE3	7:D:2002:DST:S9	2.32	0.69
8:C:355:HOH:O	1:D:161:CYS:SG	2.51	0.69
1:A:187:GLY:O	1:A:191:LEU:HG	1.92	0.68
1:D:187:GLY:O	1:D:191:LEU:HG	1.94	0.68
1:A:133:PRO:O	1:A:137:VAL:HG23	1.94	0.67
1:A:160:VAL:HA	1:A:163:GLU:OE2	1.94	0.67
2:C:124:VAL:HG22	2:C:135:ILE:HD12	1.75	0.67
1:D:138:ARG:HD2	8:D:397:HOH:O	1.94	0.66
2:B:28:VAL:O	2:B:31:PRO:HD2	1.94	0.66
2:B:2:GLN:CB	2:B:3:PRO:HD3	2.12	0.66
2:B:82:LEU:H	2:B:89:ARG:HH12	1.44	0.65
2:B:236:LYS:HG2	2:B:240:LEU:CD1	2.27	0.64
2:B:249:LYS:HZ3	2:B:253:LEU:HD13	1.63	0.64
2:C:4:TYR:CZ	2:C:8:ILE:HD11	2.32	0.64
1:D:233:LEU:HD12	1:D:237:ALA:HB2	1.79	0.64
2:C:43:ALA:O	2:C:47:LEU:HB2	1.97	0.64
2:C:227:ILE:HG12	8:C:1126:HOH:O	1.98	0.63
2:B:203:LYS:HD2	2:B:246:PHE:CE2	2.32	0.63
2:C:61:GLN:HG2	2:C:128:ARG:NH2	2.13	0.63
1:A:138:ARG:HD2	8:A:697:HOH:O	1.99	0.62
2:B:26:GLU:CB	8:B:435:HOH:O	2.39	0.62
1:D:149:SER:HB2	8:D:361:HOH:O	2.00	0.61
2:C:130:ASP:OD2	8:C:292:HOH:O	2.16	0.61
2:B:215:ARG:NH1	8:B:458:HOH:O	2.25	0.61
2:B:166:THR:OG1	2:B:171:ILE:HD11	2.01	0.61
2:B:16:LEU:HD21	2:B:67:ALA:HB1	1.83	0.61
1:D:165:MET:HE2	1:D:168:VAL:HA	1.81	0.60
2:C:197:ALA:HB3	2:C:199:GLU:OE1	2.00	0.60
2:C:172:GLU:HG3	2:C:217:MET:HE3	1.83	0.60
1:D:281:ALA:HB3	1:D:282:PRO:HD3	1.83	0.60
1:D:294:ASP:OD2	8:D:953:HOH:O	2.17	0.60
2:C:221:LYS:HD3	2:C:227:ILE:CB	2.31	0.60
2:B:168:LEU:HB3	2:B:231:ILE:HD11	1.82	0.60
2:C:215:ARG:HH11	2:C:215:ARG:HG2	1.65	0.60
2:B:103:ASN:OD1	2:B:104:VAL:HG23	2.02	0.60
2:C:125:ASP:HB3	2:C:128:ARG:HG3	1.82	0.59
2:C:92:SER:C	2:C:94:PRO:HD2	2.22	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:43:ALA:HB3	2:C:183:HIS:HE1	1.67	0.59
2:B:249:LYS:HZ2	2:B:253:LEU:HD22	1.65	0.59
1:D:61:ASP:N	1:D:61:ASP:OD2	2.35	0.59
2:C:221:LYS:CB	2:C:227:ILE:HD12	2.31	0.58
2:B:206:ASN:HB3	2:B:242:GLU:HG2	1.83	0.58
1:D:156:GLN:NE2	7:D:2002:DST:H143	2.18	0.58
2:C:4:TYR:CE2	2:C:8:ILE:HD11	2.38	0.58
2:C:43:ALA:HB3	2:C:183:HIS:CE1	2.39	0.58
1:D:254:VAL:O	1:D:257:SER:HB2	2.03	0.58
2:B:249:LYS:NZ	2:B:253:LEU:HD22	2.19	0.58
1:D:174:GLU:HG2	8:D:661:HOH:O	2.03	0.58
1:D:295:ASN:C	1:D:295:ASN:HD22	2.07	0.57
2:B:161:ILE:HA	2:B:166:THR:HG21	1.87	0.57
2:C:8:ILE:CG2	2:C:49:LEU:HD12	2.34	0.57
2:B:211:GLN:NE2	2:B:257:LEU:HB3	2.19	0.57
1:D:177:HIS:HD2	1:D:248:TYR:HE2	1.53	0.57
1:A:199:LYS:HE3	8:A:906:HOH:O	2.04	0.57
1:D:240:ASP:OD1	1:D:245:LYS:HD2	2.05	0.56
2:C:219:GLU:HG2	8:C:780:HOH:O	2.06	0.56
2:C:168:LEU:HD11	2:C:226:LEU:HG	1.86	0.56
2:B:58:ASP:H	2:B:61:GLN:NE2	1.99	0.56
2:C:221:LYS:HD3	2:C:227:ILE:HD12	1.86	0.56
1:D:185:LEU:HD13	1:D:214:GLY:HA2	1.88	0.56
2:B:131:ASP:O	2:B:135:ILE:HG13	2.05	0.55
2:B:91:VAL:HG23	8:B:882:HOH:O	2.05	0.55
1:A:258:LYS:NZ	8:A:550:HOH:O	2.27	0.55
1:A:277:PRO:O	8:A:303:HOH:O	2.18	0.55
1:A:44:LYS:HE3	8:A:715:HOH:O	2.06	0.55
1:A:47:ARG:HB2	1:A:48:PRO:HD3	1.88	0.54
2:C:91:VAL:O	2:C:91:VAL:HG23	2.07	0.54
2:C:172:GLU:HG3	2:C:217:MET:CE	2.38	0.54
1:D:30:LYS:NZ	8:D:645:HOH:O	2.40	0.54
1:D:31:ILE:O	1:D:35:MET:HG3	2.08	0.54
1:A:180:LYS:HD2	8:A:296:HOH:O	2.08	0.54
1:D:270:GLU:HG2	8:D:1011:HOH:O	2.06	0.54
2:C:108:THR:O	2:C:112:ILE:HG12	2.07	0.54
1:D:266:ARG:NH1	8:D:1051:HOH:O	2.41	0.54
1:D:267:GLU:OE1	8:D:869:HOH:O	2.18	0.54
2:B:248:GLY:O	2:B:252:GLU:HB2	2.08	0.54
2:B:82:LEU:N	2:B:89:ARG:HH12	2.05	0.53
1:D:241:LEU:HD13	8:D:1050:HOH:O	2.08	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:94:ARG:O	1:A:95:ARG:HB2	2.07	0.53
1:D:176:ILE:O	1:D:180:LYS:HB3	2.07	0.53
2:C:125:ASP:HB3	2:C:128:ARG:CG	2.39	0.53
1:D:30:LYS:HD2	1:D:108:VAL:HG21	1.90	0.53
1:D:177:HIS:HD2	1:D:248:TYR:CE2	2.25	0.53
1:D:124:VAL:HG12	1:D:124:VAL:O	2.07	0.53
1:A:138:ARG:CD	8:A:697:HOH:O	2.55	0.53
1:D:135:ARG:HG3	1:D:135:ARG:NH1	2.19	0.53
2:B:25:PRO:HB2	2:B:27:THR:CG2	2.36	0.52
2:C:218:MET:O	2:C:221:LYS:HB3	2.09	0.52
1:D:165:MET:CE	1:D:168:VAL:HA	2.38	0.52
2:C:162:VAL:O	2:C:164:GLY:N	2.42	0.52
1:A:147:ILE:O	1:A:153:VAL:HG23	2.09	0.52
1:A:186:GLN:HG3	1:A:207:ARG:HG3	1.92	0.52
2:B:229:GLU:OE1	2:B:229:GLU:HA	2.09	0.52
1:A:266:ARG:O	1:A:270:GLU:HG3	2.10	0.52
2:B:203:LYS:HD2	2:B:246:PHE:HE2	1.75	0.52
1:D:170:LEU:HD11	8:D:627:HOH:O	2.08	0.51
2:B:94:PRO:CB	8:B:968:HOH:O	2.48	0.51
2:B:200:GLU:HG3	8:B:459:HOH:O	2.10	0.51
1:A:165:MET:HG2	1:A:167:GLU:H	1.73	0.51
2:B:249:LYS:HZ3	2:B:253:LEU:CD1	2.22	0.51
2:B:90:PRO:HG3	2:B:159:GLU:HA	1.92	0.51
2:C:221:LYS:CD	2:C:227:ILE:HB	2.36	0.51
2:B:94:PRO:CG	8:B:968:HOH:O	2.59	0.51
2:B:182:MET:O	2:B:185:CYS:HB3	2.11	0.51
1:D:263:ARG:HH11	1:D:263:ARG:HB2	1.76	0.51
2:B:236:LYS:HG2	2:B:240:LEU:HD12	1.92	0.50
2:B:168:LEU:HD13	2:B:231:ILE:HD11	1.92	0.50
2:C:93:LYS:N	2:C:94:PRO:CD	2.74	0.50
2:C:97:GLN:HG3	8:C:319:HOH:O	2.12	0.50
1:D:208:LYS:HD3	1:D:271:GLN:OE1	2.12	0.50
2:B:17:LYS:HG3	2:B:37:PHE:CZ	2.47	0.50
2:C:131:ASP:N	2:C:132:PRO:HD2	2.27	0.50
2:B:54:LEU:HD11	2:B:251:ALA:HA	1.94	0.50
2:C:126:PRO:C	2:C:128:ARG:H	2.14	0.50
2:B:128:ARG:O	2:B:131:ASP:HB2	2.12	0.50
2:C:4:TYR:O	2:C:8:ILE:HG13	2.12	0.50
1:A:186:GLN:OE1	1:A:207:ARG:CD	2.60	0.49
2:C:200:GLU:OE2	2:C:246:PHE:HE1	1.95	0.49
2:C:84:LEU:HD12	2:C:89:ARG:HG2	1.93	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:7:TYR:CD2	1:A:7:TYR:C	2.85	0.49
2:B:130:ASP:O	2:B:134:ARG:HG3	2.12	0.49
1:A:86:PRO:HG2	1:A:106:GLU:OE2	2.13	0.49
2:C:163:ASP:HA	2:C:224:HIS:CE1	2.46	0.49
1:A:140:LEU:HD13	2:B:139:ILE:HG21	1.94	0.49
1:D:177:HIS:ND1	1:D:218:GLN:HG2	2.27	0.48
2:B:211:GLN:HE22	2:B:257:LEU:HB3	1.77	0.48
1:A:54:ALA:HB2	1:A:189:VAL:HG22	1.95	0.48
1:D:173:LEU:HD22	1:D:251:LEU:HD12	1.94	0.48
1:A:276:HIS:HB2	1:A:279:ARG:HD2	1.95	0.48
2:B:236:LYS:HG3	2:B:258:VAL:HB	1.95	0.48
1:D:156:GLN:HE22	7:D:2002:DST:H112	1.78	0.48
2:C:227:ILE:HG13	8:C:1126:HOH:O	2.06	0.48
2:B:162:VAL:H	2:B:166:THR:HG23	1.74	0.48
1:A:163:GLU:HB3	8:A:370:HOH:O	2.13	0.47
1:A:186:GLN:NE2	1:A:211:ASN:OD1	2.47	0.47
2:C:103:ASN:HB3	1:D:87:CYS:O	2.13	0.47
1:D:186:GLN:HE21	1:D:211:ASN:CB	2.27	0.47
1:A:168:VAL:HG13	1:A:172:HIS:CB	2.42	0.47
2:B:211:GLN:CD	2:B:257:LEU:HD23	2.35	0.47
1:A:157:VAL:O	1:A:160:VAL:HG12	2.13	0.47
2:B:227:ILE:HD12	2:B:227:ILE:N	2.29	0.47
2:B:236:LYS:HG2	2:B:240:LEU:HD11	1.95	0.47
2:B:20:ILE:O	2:B:33:HIS:HD2	1.97	0.47
1:D:170:LEU:CD2	1:D:251:LEU:HD13	2.43	0.47
2:B:1:MET:O	2:B:1:MET:HG2	2.15	0.47
1:A:30:LYS:HB3	1:A:30:LYS:HE2	1.80	0.46
1:D:243:ALA:C	1:D:245:LYS:H	2.18	0.46
2:C:227:ILE:HG22	2:C:227:ILE:O	2.14	0.46
1:A:252:ILE:O	1:A:256:LYS:HB2	2.15	0.46
2:B:210:TYR:HB2	2:B:239:ALA:HB2	1.98	0.46
2:B:215:ARG:HD3	8:B:733:HOH:O	2.15	0.46
1:D:163:GLU:HG3	1:D:245:LYS:HG2	1.97	0.46
1:D:186:GLN:HA	1:D:210:ALA:HB1	1.97	0.46
2:B:171:ILE:HG22	2:B:217:MET:CE	2.46	0.46
1:D:222:ASP:HB3	1:D:249:PRO:HD2	1.98	0.46
2:C:129:THR:O	2:C:132:PRO:HD2	2.17	0.45
2:C:204:LEU:HD23	2:C:204:LEU:HA	1.69	0.45
2:C:90:PRO:O	2:C:92:SER:N	2.49	0.45
1:D:173:LEU:CD1	1:D:246:THR:HG22	2.43	0.45
2:C:250:ASN:HA	2:C:253:LEU:CD1	2.44	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:232:ILE:O	2:C:236:LYS:HB2	2.16	0.45
1:D:9:LEU:HA	1:D:9:LEU:HD12	1.65	0.45
2:C:58:ASP:HB3	8:C:1080:HOH:O	2.15	0.45
1:A:177:HIS:HD2	1:A:248:TYR:OH	2.00	0.45
2:B:45:SER:O	2:B:48:CYS:HB3	2.17	0.45
1:A:228:LYS:O	8:A:372:HOH:O	2.21	0.45
2:C:247:HIS:CB	8:C:1129:HOH:O	2.53	0.45
2:B:167:SER:OG	1:D:33:GLU:OE2	2.18	0.45
2:B:179:TYR:CD1	2:B:216:GLY:HA3	2.52	0.45
2:B:213:THR:O	2:B:217:MET:HB2	2.16	0.44
2:C:32:MET:HG3	2:C:112:ILE:HD11	1.99	0.44
1:D:186:GLN:NE2	1:D:211:ASN:HB2	2.32	0.44
1:D:173:LEU:HD11	1:D:177:HIS:NE2	2.32	0.44
2:C:140:ILE:HD13	1:D:126:ALA:HB2	2.00	0.44
1:A:132:PRO:HA	1:A:133:PRO:HD3	1.84	0.44
2:C:221:LYS:HD3	2:C:227:ILE:CG1	2.48	0.44
2:B:221:LYS:NZ	2:B:232:ILE:HD13	2.33	0.44
1:D:97:LYS:HB3	1:D:97:LYS:HE2	1.63	0.44
1:A:208:LYS:HB3	1:A:208:LYS:HE2	1.89	0.44
2:B:240:LEU:HD22	2:B:255:SER:HB3	1.99	0.44
2:B:35:LEU:HD12	2:B:108:THR:HG21	1.99	0.44
2:C:133:ASP:OD2	1:D:133:PRO:HG2	2.18	0.44
2:C:31:PRO:HG3	2:C:100:TYR:CE1	2.52	0.44
1:D:277:PRO:O	8:D:1053:HOH:O	2.21	0.44
1:A:71:ALA:HB2	1:A:124:VAL:HG23	2.00	0.43
2:C:93:LYS:O	2:C:94:PRO:C	2.57	0.43
2:B:221:LYS:CG	2:B:227:ILE:HG12	2.34	0.43
2:B:30:GLY:N	2:B:31:PRO:CD	2.81	0.43
2:B:82:LEU:O	2:B:84:LEU:HG	2.17	0.43
2:B:90:PRO:CD	2:B:159:GLU:HB3	2.48	0.43
1:D:182:ALA:O	1:D:186:GLN:HG3	2.18	0.43
2:C:26:GLU:HG2	8:C:803:HOH:O	2.17	0.43
1:D:255:GLU:HA	1:D:255:GLU:OE1	2.19	0.43
1:A:96:GLY:C	1:A:97:LYS:HD2	2.39	0.43
2:B:23:ARG:HB2	2:B:24:PRO:CD	2.40	0.43
2:B:128:ARG:NH1	2:B:128:ARG:CG	2.67	0.43
1:D:222:ASP:HB3	1:D:249:PRO:CD	2.49	0.43
2:B:41:ALA:HB1	8:B:316:HOH:O	2.18	0.43
1:D:68:ALA:HA	1:D:124:VAL:CG2	2.45	0.43
2:B:236:LYS:O	2:B:240:LEU:HG	2.18	0.43
1:D:292:TYR:HE2	8:D:381:HOH:O	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:1:MET:HG3	2:B:256:SER:HB3	2.01	0.42
2:C:85:THR:OG1	2:C:86:ASP:N	2.52	0.42
1:D:124:VAL:O	1:D:124:VAL:CG1	2.67	0.42
1:A:290:ILE:CD1	3:A:2003:IPE:H51	2.49	0.42
2:B:240:LEU:CD2	2:B:255:SER:HB3	2.49	0.42
1:A:199:LYS:N	1:A:202:GLU:OE1	2.47	0.42
2:B:237:GLU:O	2:B:241:GLU:HG3	2.19	0.42
2:C:138:VAL:HG12	2:C:142:ILE:HD12	2.01	0.42
2:C:249:LYS:O	2:C:253:LEU:HG	2.19	0.42
2:B:71:LEU:CD2	2:B:112:ILE:HG23	2.44	0.42
2:C:162:VAL:C	2:C:164:GLY:N	2.73	0.42
1:A:167:GLU:HG2	1:A:167:GLU:O	2.20	0.42
1:D:71:ALA:HB1	1:D:120:ALA:HB1	2.02	0.42
2:B:171:ILE:HG22	2:B:217:MET:HE3	2.01	0.42
1:D:97:LYS:HB2	1:D:98:PRO:CD	2.49	0.42
1:D:139:VAL:HG21	1:D:195:LEU:HG	2.00	0.42
1:D:233:LEU:HD12	1:D:237:ALA:CB	2.49	0.42
1:A:281:ALA:HB3	1:A:282:PRO:HD3	2.02	0.42
2:B:8:ILE:HG23	2:B:49:LEU:HD12	2.01	0.42
1:A:129:LYS:CE	8:A:415:HOH:O	2.68	0.42
2:B:247:HIS:C	2:B:251:ALA:HB3	2.41	0.41
1:D:139:VAL:HG13	1:D:191:LEU:HD22	2.01	0.41
1:D:186:GLN:HA	1:D:210:ALA:CB	2.50	0.41
2:C:251:ALA:O	2:C:255:SER:HB2	2.18	0.41
2:C:1:MET:N	8:C:920:HOH:O	2.52	0.41
2:C:112:ILE:HD13	2:C:112:ILE:HA	1.78	0.41
2:B:161:ILE:HD13	2:B:171:ILE:HD12	2.02	0.41
2:C:14:ARG:NE	8:C:922:HOH:O	2.42	0.41
2:B:158:GLU:OE1	2:B:178:LYS:NZ	2.49	0.41
1:A:225:ASP:O	1:A:228:LYS:HD2	2.21	0.41
1:A:206:LEU:HD23	1:A:206:LEU:HA	1.85	0.41
2:C:88:SER:OG	1:D:107:SER:HB2	2.20	0.41
2:B:232:ILE:H	2:B:232:ILE:HG12	1.66	0.41
2:C:217:MET:O	2:C:227:ILE:HD11	2.21	0.41
2:C:1:MET:HE2	8:C:919:HOH:O	2.21	0.41
1:A:19:LEU:HD23	1:A:19:LEU:HA	1.85	0.41
2:C:206:ASN:O	2:C:210:TYR:CD1	2.73	0.41
1:A:8:MET:HE3	8:A:317:HOH:O	2.19	0.41
2:C:65:ALA:O	2:C:69:ILE:HG13	2.21	0.41
1:A:168:VAL:HG13	1:A:172:HIS:CG	2.56	0.40
1:A:139:VAL:HG13	1:A:191:LEU:HD22	2.03	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:170:LEU:CD1	8:D:627:HOH:O	2.69	0.40
1:D:225:ASP:OD2	1:D:240:ASP:HB2	2.22	0.40
2:C:32:MET:HG3	2:C:112:ILE:CD1	2.52	0.40
2:C:139:ILE:HG21	1:D:140:LEU:HD13	2.03	0.40
1:A:186:GLN:OE1	1:A:207:ARG:HD3	2.21	0.40
2:B:165:ASN:O	2:B:165:ASN:CG	2.60	0.40
2:B:93:LYS:N	2:B:94:PRO:HD2	2.36	0.40
2:C:162:VAL:H	2:C:166:THR:HG21	1.87	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
8:A:843:HOH:O	8:C:805:HOH:O[4_555]	1.99	0.21
8:B:589:HOH:O	8:D:588:HOH:O[3_555]	2.07	0.13

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	275/295 (93%)	263 (96%)	12 (4%)	0	100	100
1	D	293/295 (99%)	280 (96%)	12 (4%)	1 (0%)	46	35
2	B	257/274 (94%)	237 (92%)	14 (5%)	6 (2%)	8	1
2	C	257/274 (94%)	230 (90%)	18 (7%)	9 (4%)	4	0
All	All	1082/1138 (95%)	1010 (93%)	56 (5%)	16 (2%)	13	3

All (16) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	227	ILE

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Mol	Chain	Res	Type
2	C	91	VAL
2	C	163	ASP
2	C	227	ILE
2	B	2	GLN
2	B	163	ASP
2	B	225	GLN
2	C	86	ASP
2	C	164	GLY
1	D	42	GLY
2	C	94	PRO
2	C	127	ALA
2	C	225	GLN
2	C	226	LEU
2	B	226	LEU
2	B	244	GLY

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	222/234 (95%)	218 (98%)	4 (2%)	66	60
1	D	234/234 (100%)	231 (99%)	3 (1%)	76	72
2	B	201/214 (94%)	193 (96%)	8 (4%)	38	23
2	C	201/214 (94%)	197 (98%)	4 (2%)	63	55
All	All	858/896 (96%)	839 (98%)	19 (2%)	60	51

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

Mol	Chain	Res	Type
1	A	97	LYS
1	A	135	ARG
1	A	160	VAL
1	A	189	VAL
2	B	45	SER
2	B	85	THR

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Mol	Chain	Res	Type
2	B	128	ARG
2	B	134	ARG
2	B	227	ILE
2	B	229	GLU
2	B	232	ILE
2	B	255	SER
2	C	58	ASP
2	C	59	ARG
2	C	131	ASP
2	C	199	GLU
1	D	174	GLU
1	D	199	LYS
1	D	295	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	177	HIS
2	B	33	HIS
2	B	61	GLN
2	C	183	HIS
2	C	225	GLN
2	C	230	ASN
1	D	156	GLN
1	D	186	GLN
1	D	211	ASN
1	D	295	ASN

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

Of 9 ligands modelled in this entry, 4 are monoatomic - leaving 5 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
3	IPE	A	2003	-	10,13,13	2.45	2 (20%)	14,19,19	2.71	6 (42%)
4	PPV	A	2004	5	6,8,8	1.39	0	11,13,13	2.07	1 (9%)
6	EDO	C	2005	-	3,3,3	0.71	0	2,2,2	0.52	0
3	IPE	D	2001	-	10,13,13	2.19	1 (10%)	14,19,19	2.28	3 (21%)
7	DST	D	2002	5	8,13,13	1.30	1 (12%)	11,19,19	2.68	4 (36%)

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
3	IPE	A	2003	-	-	0/13/13/13	0/0/0/0
4	PPV	A	2004	5	-	0/6/6/6	0/0/0/0
6	EDO	C	2005	-	-	0/1/1/1	0/0/0/0
3	IPE	D	2001	-	-	0/13/13/13	0/0/0/0
7	DST	D	2002	5	-	0/9/13/13	0/0/0/0

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	D	2002	DST	C11-C12	-2.15	1.35	1.51
3	A	2003	IPE	PA-O1A	2.24	1.59	1.51
3	D	2001	IPE	C4-C3	6.46	1.54	1.33
3	A	2003	IPE	C4-C3	6.98	1.55	1.33

All (14) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	A	2004	PPV	P2-OPP-P1	-6.49	110.91	132.67

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	D	2002	DST	O4-P1-O5	-3.99	97.75	110.58
3	A	2003	IPE	PA-O3A-PB	-3.93	119.49	132.67
3	D	2001	IPE	PA-O3A-PB	-2.60	123.95	132.67
3	A	2003	IPE	O3B-PB-O1B	-2.49	97.90	107.38
7	D	2002	DST	C13-C12-C11	2.01	125.96	111.38
3	A	2003	IPE	O1-C1-C2	2.04	118.42	108.63
3	A	2003	IPE	O3B-PB-O3A	2.26	115.33	105.09
3	D	2001	IPE	O1B-PB-O3A	2.49	116.41	105.09
3	A	2003	IPE	O1B-PB-O3A	3.04	118.88	105.09
7	D	2002	DST	O4-P1-O2	4.07	123.54	105.09
7	D	2002	DST	O8-P3-O7	5.73	124.36	110.05
3	D	2001	IPE	O3A-PA-O1	6.42	119.97	102.94
3	A	2003	IPE	O3A-PA-O1	7.41	122.60	102.94

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	2003	IPE	1	0
7	D	2002	DST	3	0

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	279/295 (94%)	0.32	21 (7%)	17 26	27, 39, 68, 99	0
1	D	295/295 (100%)	0.49	25 (8%)	13 21	26, 41, 91, 109	0
2	B	259/274 (94%)	0.95	48 (18%)	2 2	29, 46, 116, 126	0
2	C	259/274 (94%)	0.86	39 (15%)	3 5	27, 46, 96, 115	0
All	All	1092/1138 (95%)	0.64	133 (12%)	5 9	26, 42, 98, 126	0

All (133) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	251	LEU	12.7
2	B	248	GLY	10.6
2	B	257	LEU	9.4
2	B	246	PHE	8.9
2	B	250	ASN	8.4
2	B	249	LYS	8.2
2	C	129	THR	8.2
1	A	166	ALA	7.6
2	C	91	VAL	7.5
2	B	1	MET	7.4
2	B	228	ASP	7.1
2	C	127	ALA	7.1
2	C	126	PRO	7.0
2	B	226	LEU	7.0
1	A	164	GLY	6.8
1	A	165	MET	6.8
1	D	168	VAL	6.7
1	D	243	ALA	6.6
1	D	165	MET	6.6
1	D	244	ASP	6.4
2	B	93	LYS	6.2

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Mol	Chain	Res	Type	RSRZ
2	C	93	LYS	6.2
1	D	242	VAL	6.0
1	D	167	GLU	6.0
1	D	166	ALA	5.9
2	B	225	GLN	5.9
2	B	251	ALA	5.8
2	B	3	PRO	5.7
2	C	249	LYS	5.7
2	C	225	GLN	5.4
2	B	222	ASN	5.1
2	B	221	LYS	5.1
2	C	226	LEU	5.0
2	C	222	ASN	4.9
1	D	246	THR	4.8
2	C	128	ARG	4.8
2	C	259	ALA	4.7
2	B	223	SER	4.7
1	D	252	ILE	4.6
2	B	96	ILE	4.6
2	B	238	LEU	4.5
1	D	164	GLY	4.5
2	B	256	SER	4.5
2	B	231	ILE	4.3
2	C	94	PRO	4.2
2	B	258	VAL	4.1
2	B	254	MET	4.1
2	C	124	VAL	4.1
1	D	170	LEU	4.1
1	D	171	ASP	4.0
2	B	224	HIS	3.9
1	D	254	VAL	3.8
2	C	162	VAL	3.8
2	B	90	PRO	3.8
2	C	125	ASP	3.7
1	A	167	GLU	3.7
2	C	224	HIS	3.7
1	D	245	LYS	3.7
2	C	223	SER	3.6
2	C	130	ASP	3.6
2	B	2	GLN	3.6
2	B	24	PRO	3.6
2	C	221	LYS	3.5

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Mol	Chain	Res	Type	RSRZ
1	A	161	CYS	3.4
1	A	276	HIS	3.4
2	C	95	ALA	3.3
2	C	163	ASP	3.3
2	B	162	VAL	3.2
2	C	230	ASN	3.2
1	D	184	LEU	3.2
2	C	92	SER	3.1
1	A	278	HIS	3.1
2	B	227	ILE	3.1
2	C	227	ILE	3.1
2	B	245	GLY	3.1
2	B	91	VAL	3.0
2	B	164	GLY	3.0
1	A	96	GLY	3.0
2	C	244	GLY	3.0
2	C	123	SER	2.9
2	B	165	ASN	2.9
2	C	247	HIS	2.9
1	D	239	LYS	2.8
1	A	72	VAL	2.8
1	A	184	LEU	2.8
2	C	164	GLY	2.8
2	B	243	LEU	2.8
1	A	228	LYS	2.7
1	A	185	LEU	2.7
2	B	95	ALA	2.7
1	A	295	ASN	2.7
2	B	229	GLU	2.7
2	B	235	LEU	2.7
2	B	237	GLU	2.7
2	C	253	LEU	2.6
1	D	185	LEU	2.6
2	B	240	LEU	2.6
2	B	244	GLY	2.6
1	D	259	GLU	2.6
1	A	95	ARG	2.5
1	A	117	LEU	2.5
2	B	27	THR	2.5
1	A	147	ILE	2.5
2	B	92	SER	2.4
2	B	255	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	A	75	ILE	2.4
1	D	147	ILE	2.4
2	B	230	ASN	2.4
2	C	182	MET	2.4
2	B	259	ALA	2.4
2	B	253	LEU	2.3
1	A	163	GLU	2.3
1	D	162	SER	2.3
1	D	238	GLY	2.3
2	C	186	GLY	2.3
2	C	165	ASN	2.3
2	B	241	GLU	2.2
2	C	89	ARG	2.2
2	C	90	PRO	2.2
1	A	162	SER	2.2
2	B	252	GLU	2.2
1	A	274	HIS	2.2
2	C	258	VAL	2.2
2	C	199	GLU	2.2
2	C	122	GLY	2.1
2	B	247	HIS	2.1
2	B	51	ALA	2.1
1	D	161	CYS	2.1
1	D	163	GLU	2.0
1	D	4	PHE	2.0
2	C	142	ILE	2.0
2	C	47	LEU	2.0
1	A	213	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 carbohydrates in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron

density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors(Å ²)	Q<0.9
6	EDO	C	2005	4/4	0.96	0.19	2.43	36,37,37,38	0
3	IPE	A	2003	14/14	0.89	0.20	1.15	72,81,88,88	0
7	DST	D	2002	14/14	0.96	0.15	-0.08	49,55,62,62	0
3	IPE	D	2001	14/14	0.98	0.12	-0.21	39,40,45,48	0
5	MG	D	3002	1/1	0.95	0.06	-0.86	49,49,49,49	0
4	PPV	A	2004	9/9	0.93	0.10	-1.09	76,76,77,78	0
5	MG	D	3001	1/1	0.93	0.06	-1.16	53,53,53,53	0
5	MG	A	3003	1/1	0.94	0.11	-	50,50,50,50	0
5	MG	A	3004	1/1	0.83	0.07	-	58,58,58,58	0

6.5 Other polymers

There are no such residues in this entry.