



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 5, 2026 – 05:13 PM UTC

PDB ID : 1H18 / pdb_00001h18
Title : Pyruvate Formate-Lyase (E.coli) in complex with Pyruvate
Authors : Becker, A.; Kabsch, W.
Deposited on : 2002-07-04
Resolution : 2.30 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity : 4-5-2 with Phenix2.0
Mogul : 2022.3.0, CSD as543be (2022)
Xtriage (Phenix) : 2.0
EDS : 3.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4 : 9.0.010 (Gargrove)
Density-Fitness : 1.0.12
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

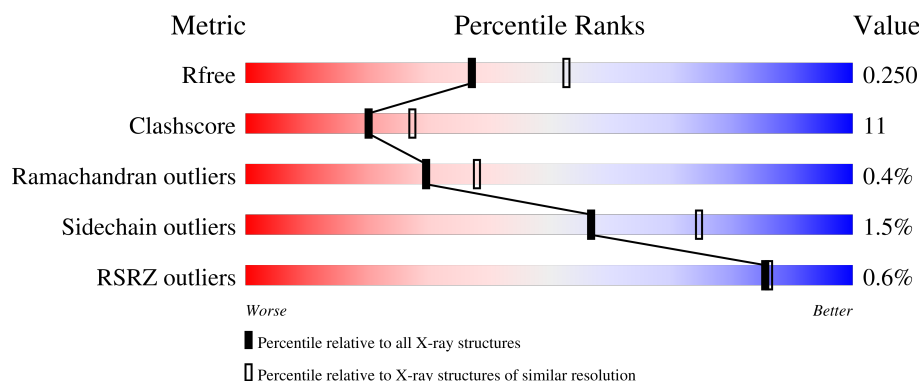
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	180053	6319 (2.30-2.30)
Clashscore	190562	6919 (2.30-2.30)
Ramachandran outliers	187476	6854 (2.30-2.30)
Sidechain outliers	187428	6854 (2.30-2.30)
RSRZ outliers	180081	6325 (2.30-2.30)

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 $\leq 5\%$. 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	759	77% 22% .
1	B	759	72% 26% .

2 Entry composition [i](#)

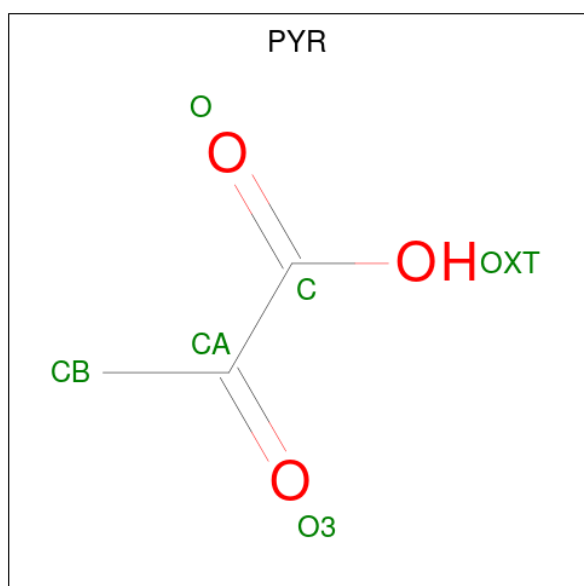
There are 6 unique types of molecules in this entry. The entry contains 13667 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 FORMATE ACETYLTRANSFERASE 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	759	Total	C	N	O	S	0	39	0
			6172	3906	1050	1174	42			
1	B	759	Total	C	N	O	S	0	39	0
			6172	3906	1050	1174	42			

- Molecule 2 is PYRUVIC ACID (CCD ID: PYR) (formula: $C_3H_4O_3$).

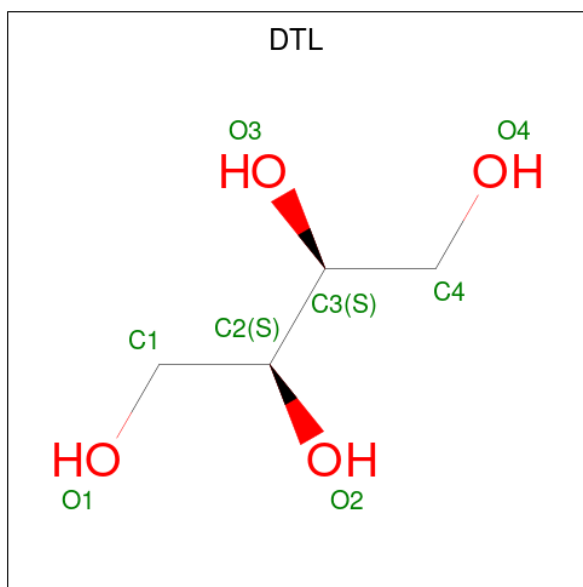


Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			6	3	3		
2	B	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is SODIUM ION (CCD ID: NA) (formula: Na).

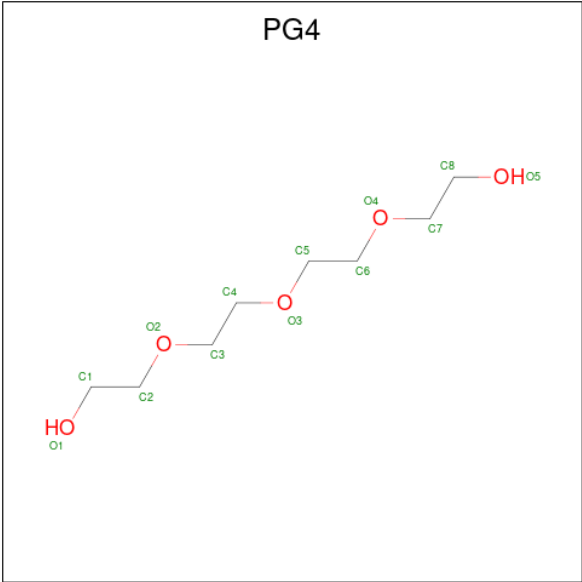
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	3	Total	Na	0	0
			3	3		
3	B	4	Total	Na	0	0
			4	4		

- Molecule 4 is L-TREITOL (CCD ID: DTL) (formula: $C_4H_{10}O_4$).



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

- Molecule 5 is TETRAETHYLENE GLYCOL (CCD ID: PG4) (formula: $C_8H_{18}O_5$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
5	A	1	Total	C	O	0	0
			13	8	5		
5	A	1	Total	C	O	0	0
			13	8	5		
5	A	1	Total	C	O	0	0
			13	8	5		
5	B	1	Total	C	O	0	0
			13	8	5		
5	B	1	Total	C	O	0	0
			13	8	5		

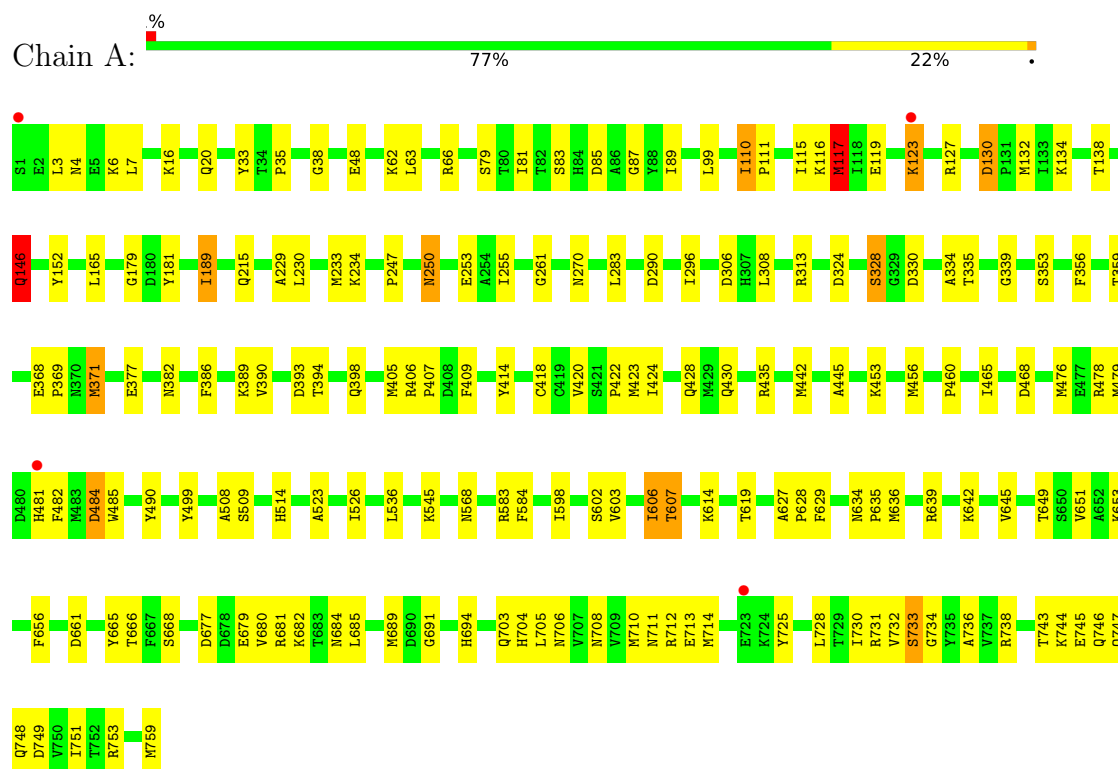
- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	633	Total	O	0	0
			633	633		
6	B	574	Total	O	0	0
			574	574		

3 Residue-property plots

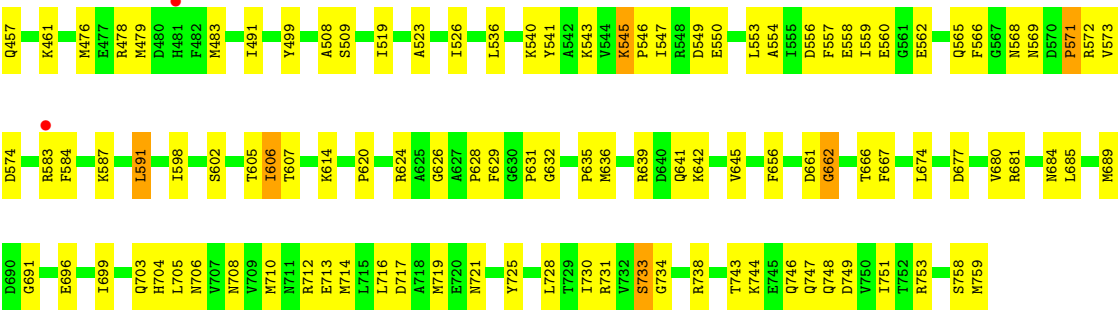
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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: FORMATE ACETYLTRANSFERASE 1



• Molecule 1: FORMATE ACETYLTRANSFERASE 1





4 Data and refinement statistics

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	158.91Å 158.91Å 159.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	15.00 – 2.30 15.00 – 2.30	Depositor EDS
% Data completeness (in resolution range)	95.2 (15.00-2.30) 94.9 (15.00-2.30)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	3.55 (at 2.29Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.183 , 0.234 0.198 , 0.250	Depositor DCC
R_{free} test set	1734 reflections (2.00%)	wwPDB-VP
Wilson B-factor (Å ²)	47.5	Xtriage
Anisotropy	0.076	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 48.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.016 for -h,l,k 0.004 for -l,-k,-h	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	13667	wwPDB-VP
Average B, all atoms (Å ²)	52.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.27% 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: NA, PG4, PYR, DTL

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	2.42	3/6477 (0.0%)	1.29	39/8742 (0.4%)
1	B	0.55	0/6477	0.99	34/8742 (0.4%)
All	All	1.75	3/12954 (0.0%)	1.15	73/17484 (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	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	146[A]	GLN	CD-NE2	133.85	4.14	1.33
1	A	146[B]	GLN	CD-NE2	133.85	4.14	1.33
1	A	117	MET	SD-CE	-5.07	1.66	1.79

All (73) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	146[A]	GLN	CG-CD-NE2	-47.83	44.66	116.40
1	A	146[B]	GLN	CG-CD-NE2	-47.83	44.66	116.40
1	A	146[A]	GLN	OE1-CD-NE2	23.82	146.42	122.60
1	A	146[B]	GLN	OE1-CD-NE2	23.82	146.42	122.60
1	B	89	ILE	N-CA-C	-11.32	100.92	111.45
1	A	146[A]	GLN	CA-CB-CG	9.59	133.28	114.10
1	A	146[B]	GLN	CA-CB-CG	9.59	133.28	114.10
1	A	89	ILE	N-CA-C	-9.58	102.54	111.45

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	662	GLY	N-CA-C	9.31	120.79	111.95
1	A	189	ILE	N-CA-C	7.43	117.55	110.42
1	B	425	VAL	N-CA-C	7.35	118.85	108.93
1	B	110[A]	ILE	CA-C-N	7.06	126.67	119.19
1	B	110[A]	ILE	C-N-CA	7.06	126.67	119.19
1	B	110[B]	ILE	CA-C-N	7.06	126.67	119.19
1	B	110[B]	ILE	C-N-CA	7.06	126.67	119.19
1	A	111	PRO	N-CA-C	7.04	122.66	113.86
1	A	146[A]	GLN	CB-CG-CD	6.82	124.20	112.60
1	A	146[B]	GLN	CB-CG-CD	6.82	124.20	112.60
1	B	410	ASN	N-CA-C	-6.76	102.06	111.52
1	A	394	THR	N-CA-C	6.68	122.93	114.31
1	B	189	ILE	N-CA-C	6.56	116.72	110.42
1	A	508	ALA	N-CA-C	6.37	118.75	111.11
1	A	635	PRO	N-CA-C	-6.28	101.99	111.41
1	A	83	SER	N-CA-C	6.17	117.67	111.07
1	B	43	LEU	N-CA-C	6.12	118.86	110.24
1	B	353	SER	N-CA-C	-6.02	104.72	111.28
1	A	110[A]	ILE	CA-C-N	6.01	125.63	119.56
1	A	110[A]	ILE	C-N-CA	6.01	125.63	119.56
1	A	110[B]	ILE	CA-C-N	6.01	125.63	119.56
1	A	110[B]	ILE	C-N-CA	6.01	125.63	119.56
1	A	405	MET	N-CA-C	5.96	119.02	111.69
1	B	323	TYR	N-CA-C	-5.95	104.70	111.07
1	B	641[A]	GLN	N-CA-C	5.95	119.59	112.93
1	B	641[B]	GLN	N-CA-C	5.95	119.59	112.93
1	B	179	GLY	N-CA-C	-5.94	104.48	112.14
1	B	431	PHE	N-CA-C	-5.87	97.65	107.99
1	B	499	TYR	N-CA-C	-5.86	104.81	111.14
1	B	508	ALA	N-CA-C	5.80	118.07	111.11
1	A	328	SER	N-CA-C	5.76	118.14	110.53
1	A	484	ASP	N-CA-C	-5.73	104.95	111.14
1	A	130[A]	ASP	CA-C-N	5.67	125.78	119.32
1	A	130[A]	ASP	C-N-CA	5.67	125.78	119.32
1	A	130[B]	ASP	CA-C-N	5.67	125.78	119.32
1	A	130[B]	ASP	C-N-CA	5.67	125.78	119.32
1	A	308	LEU	N-CA-C	-5.65	105.04	111.14
1	B	213	LEU	N-CA-C	5.64	117.87	111.11
1	A	634	ASN	N-CA-C	5.59	117.79	110.08
1	B	591	LEU	N-CA-C	5.54	117.88	110.35
1	A	334	ALA	N-CA-C	-5.53	100.60	109.39
1	A	499	TYR	N-CA-C	-5.51	105.28	111.28

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	629	PHE	N-CA-C	-5.43	102.72	110.59
1	A	619	THR	CA-C-N	5.42	125.76	119.47
1	A	619	THR	C-N-CA	5.42	125.76	119.47
1	B	204	GLN	N-CA-C	5.42	117.18	111.28
1	A	335	THR	N-CA-C	5.38	118.03	109.96
1	B	738	ARG	N-CA-C	-5.35	101.06	109.25
1	A	629	PHE	N-CA-C	-5.35	102.97	110.35
1	A	523	ALA	N-CA-C	5.28	117.36	108.96
1	B	242[A]	TYR	N-CA-C	5.22	117.99	109.59
1	B	242[B]	TYR	N-CA-C	5.22	117.99	109.59
1	B	137	PHE	N-CA-C	5.21	118.76	112.93
1	B	523	ALA	N-CA-C	5.15	116.56	109.15
1	B	130[A]	ASP	CA-C-N	5.10	125.13	119.32
1	B	130[A]	ASP	C-N-CA	5.10	125.13	119.32
1	B	130[B]	ASP	CA-C-N	5.10	125.13	119.32
1	B	130[B]	ASP	C-N-CA	5.10	125.13	119.32
1	A	738	ARG	N-CA-C	-5.09	101.46	109.25
1	A	353	SER	N-CA-C	-5.08	105.64	111.07
1	B	334	ALA	N-CA-C	-5.06	101.34	109.39
1	B	328	SER	N-CA-C	5.02	117.33	110.55
1	B	483	MET	N-CA-C	-5.02	105.91	112.23
1	A	339	GLY	N-CA-C	5.00	119.45	110.95
1	A	732	VAL	N-CA-C	5.00	119.75	109.34

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	146[A]	GLN	Sidechain
1	A	181	TYR	Sidechain

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	6172	0	6153	133	0
1	B	6172	0	6153	149	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	6	0	0	2	0
2	B	6	0	0	3	0
3	A	3	0	0	0	0
3	B	4	0	0	0	0
4	A	16	0	20	2	0
4	B	16	0	20	2	0
5	A	39	0	54	2	0
5	B	26	0	36	1	0
6	A	633	0	0	8	0
6	B	574	0	0	14	0
All	All	13667	0	12436	282	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 11.

All (282) 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:146[A]:GLN:CB	1:A:146[A]:GLN:CG	1.87	1.49
1:A:132:MET:SD	6:A:2183:HOH:O	1.97	1.22
1:A:117:MET:HE3	1:A:117:MET:H	1.20	1.03
1:A:110[B]:ILE:HD12	1:A:270:ASN:HB3	1.49	0.94
1:A:146[A]:GLN:CG	1:A:146[A]:GLN:CA	2.46	0.93
1:B:606:ILE:HG22	1:B:607:THR:H	1.32	0.92
1:A:117:MET:HE3	1:A:117:MET:N	1.86	0.89
1:A:602:SER:HB3	1:A:661:ASP:HB3	1.54	0.88
1:B:252:GLN:HG2	6:B:2305:HOH:O	1.75	0.85
1:B:110[B]:ILE:HD12	1:B:270:ASN:HB3	1.58	0.84
1:B:442:MET:HE1	1:B:536:LEU:HG	1.60	0.82
1:B:543:LYS:HB2	1:B:560:GLU:HB3	1.60	0.81
1:A:418:CYS:SG	2:A:1001:PYR:CA	2.70	0.80
1:A:748[B]:GLN:NE2	1:A:751[B]:ILE:HD11	1.97	0.78
1:B:250:ASN:HD21	1:B:253:GLU:HG3	1.46	0.78
1:B:685:LEU:HG	1:B:689:MET:HE2	1.66	0.78
1:B:79:SER:HB3	1:B:110[A]:ILE:HD13	1.66	0.75
1:B:418:CYS:SG	2:B:1001:PYR:CA	2.74	0.75
1:A:649:THR:O	1:A:653:LYS:HG3	1.89	0.73
1:A:250:ASN:ND2	1:A:253:GLU:H	1.85	0.73
1:B:250:ASN:ND2	1:B:253:GLU:H	1.87	0.72
1:A:250:ASN:C	1:A:250:ASN:HD22	1.98	0.71
1:B:476:MET:HE3	1:B:479[B]:MET:HB3	1.73	0.70

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:526[A]:ILE:HD11	1:A:603:VAL:HG22	1.72	0.70
1:A:16:LYS:HG2	1:A:20[A]:GLN:OE1	1.92	0.69
1:A:710:MET:HE1	1:A:730[B]:ILE:HG22	1.73	0.69
1:A:706:ASN:HD21	1:A:734:GLY:N	1.90	0.69
1:B:568:ASN:HB3	1:B:642[B]:LYS:HD2	1.74	0.69
1:A:16:LYS:HG2	1:A:20[B]:GLN:HE22	1.59	0.68
1:B:476:MET:HE3	1:B:479[A]:MET:HB3	1.76	0.66
1:B:606:ILE:HG22	1:B:607:THR:N	2.08	0.66
1:B:442:MET:HA	1:B:479[A]:MET:SD	2.36	0.66
1:A:666:THR:HA	1:A:706:ASN:HB2	1.78	0.66
1:B:250:ASN:ND2	1:B:253:GLU:HG3	2.11	0.66
1:B:553:LEU:HD22	6:B:2427:HOH:O	1.96	0.65
1:B:545:LYS:HD2	1:B:558:GLU:OE1	1.96	0.65
1:A:146[A]:GLN:CD	1:A:146[A]:GLN:HA	2.22	0.64
1:B:602:SER:HB3	1:B:661:ASP:HB3	1.80	0.64
1:A:3:LEU:HD12	1:A:3:LEU:O	1.97	0.64
1:A:418:CYS:SG	2:A:1001:PYR:C	2.86	0.64
1:A:568:ASN:HB3	1:A:642[B]:LYS:HD2	1.80	0.64
1:B:706:ASN:HD21	1:B:734:GLY:N	1.97	0.63
1:B:545:LYS:HB3	1:B:558:GLU:HG2	1.79	0.63
1:A:115:ILE:HG21	1:A:138[B]:THR:HG22	1.80	0.63
1:B:330:ASP:HB2	4:B:9010:DTL:H4C2	1.81	0.63
1:A:526[B]:ILE:HD11	1:A:584:PHE:CD2	2.34	0.63
1:B:250:ASN:C	1:B:250:ASN:HD22	2.07	0.63
1:A:250:ASN:HD21	1:A:253:GLU:H	1.47	0.62
1:B:442:MET:CE	1:B:536:LEU:HG	2.28	0.62
1:A:146[A]:GLN:CG	1:A:146[A]:GLN:HA	2.27	0.62
1:B:428[A]:GLN:HG2	1:B:519:ILE:HB	1.83	0.61
1:A:116:LYS:HB2	1:A:117:MET:CE	2.31	0.61
1:A:442:MET:HE2	1:A:479[B]:MET:SD	2.41	0.60
1:A:743:THR:O	1:A:747:GLN:HG3	2.00	0.60
1:B:295[A]:LYS:HD2	6:B:2145:HOH:O	2.01	0.60
1:B:143:THR:OG1	1:B:146[A]:GLN:HG2	2.01	0.60
1:B:583:ARG:O	1:B:587:LYS:HG3	2.02	0.60
1:A:406:ARG:HB3	1:A:407:PRO:HD3	1.84	0.60
1:B:117:MET:HE3	1:B:117:MET:N	2.17	0.60
1:B:714[A]:MET:HE1	1:B:728:LEU:HD11	1.83	0.60
1:A:117:MET:H	1:A:117:MET:CE	2.06	0.60
1:A:465:ILE:HD11	1:A:478:ARG:HG3	1.84	0.59
1:B:680:VAL:HG12	1:B:684:ASN:ND2	2.17	0.59
1:B:3:LEU:HD12	1:B:3:LEU:O	2.03	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:48:GLU:HB3	6:A:2071:HOH:O	2.02	0.59
1:B:233[A]:MET:HE3	1:B:261:GLY:HA2	1.83	0.59
1:B:666:THR:HA	1:B:706:ASN:HB2	1.84	0.59
1:B:748[B]:GLN:NE2	1:B:751[B]:ILE:HD11	2.18	0.58
1:A:4:ASN:OD1	1:A:6:LYS:HB2	2.03	0.58
1:B:568:ASN:HB3	1:B:642[B]:LYS:CD	2.33	0.58
1:A:614:LYS:HD2	1:A:745:GLU:OE1	2.04	0.58
1:B:313:ARG:HG2	1:B:368:GLU:O	2.04	0.58
1:B:725:TYR:HB3	1:B:728:LEU:HB2	1.84	0.58
1:B:566:PHE:O	1:B:635:PRO:HB3	2.03	0.58
1:B:710:MET:SD	6:B:2523:HOH:O	2.57	0.58
1:A:290:ASP:HB3	1:A:296:ILE:HG12	1.86	0.57
1:B:491:ILE:HG13	1:B:591:LEU:HD12	1.86	0.57
1:B:545:LYS:NZ	1:B:545:LYS:HB2	2.19	0.57
1:B:146[A]:GLN:NE2	6:B:2190:HOH:O	2.37	0.57
1:A:748[B]:GLN:HE22	1:A:751[B]:ILE:HD11	1.69	0.56
1:A:636:MET:HB2	1:A:639:ARG:HD2	1.88	0.56
1:A:233[A]:MET:CE	1:A:261:GLY:HA2	2.36	0.56
1:A:16:LYS:HG2	1:A:20[B]:GLN:NE2	2.21	0.55
1:A:115:ILE:HG22	6:A:2188:HOH:O	2.05	0.55
1:B:4:ASN:OD1	1:B:6:LYS:HB2	2.07	0.55
1:B:406:ARG:HB3	1:B:407:PRO:HD3	1.89	0.55
1:B:541:TYR:CE2	1:B:572:ARG:HD2	2.41	0.55
1:B:60:GLY:O	1:B:64:GLU:HG3	2.07	0.55
1:B:398:GLN:HB3	1:B:731:ARG:NH2	2.22	0.55
1:B:744:LYS:O	1:B:748[A]:GLN:HG2	2.07	0.55
1:B:614:LYS:HA	1:B:626:GLY:HA2	1.88	0.55
1:A:306:ASP:OD1	1:A:359:THR:HG22	2.07	0.55
1:A:706:ASN:HD21	1:A:734:GLY:H	1.54	0.54
1:A:127:ARG:NH1	6:A:2175:HOH:O	2.27	0.54
1:A:442:MET:HE2	1:A:479[A]:MET:HE2	1.89	0.54
1:B:565:GLN:C	1:B:573:VAL:HG11	2.33	0.54
1:A:115:ILE:HG21	1:A:138[B]:THR:CG2	2.38	0.54
1:A:691:GLY:HA3	5:A:9013:PG4:H61	1.90	0.54
1:A:602:SER:HB3	1:A:661:ASP:CB	2.32	0.53
1:B:719:MET:SD	1:B:751[B]:ILE:HG21	2.48	0.53
1:B:457:GLN:OE1	1:B:461:LYS:HA	2.08	0.53
1:A:330:ASP:HB2	4:A:9010:DTL:H4C2	1.89	0.53
1:B:134:LYS:O	1:B:138[B]:THR:HG23	2.09	0.53
1:B:37:GLU:HA	6:B:2053:HOH:O	2.08	0.53
1:B:545:LYS:HB3	1:B:558:GLU:CG	2.39	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:85:ASP:HB3	6:A:2142:HOH:O	2.09	0.53
1:B:73:PHE:CZ	1:B:127:ARG:HG3	2.44	0.52
1:A:476:MET:HE2	1:A:583:ARG:CZ	2.38	0.52
1:A:406:ARG:HG2	1:A:406:ARG:HH11	1.75	0.52
1:B:526[B]:ILE:HD11	1:B:584:PHE:CE2	2.44	0.52
1:A:744:LYS:O	1:A:748[A]:GLN:HG2	2.10	0.52
1:B:656:PHE:CD1	1:B:703:GLN:HG3	2.44	0.52
1:B:691:GLY:HA3	5:B:9013:PG4:H61	1.92	0.52
1:A:130[A]:ASP:OD2	1:A:132:MET:HB3	2.10	0.52
1:B:284:ASP:HB2	1:B:352:ASN:HB2	1.92	0.52
1:B:549:ASP:OD2	1:B:553:LEU:HB3	2.09	0.52
1:A:313:ARG:HG2	1:A:368:GLU:O	2.09	0.52
1:A:430:GLN:HG3	6:A:2492:HOH:O	2.10	0.52
1:B:677:ASP:OD2	1:B:680:VAL:HG23	2.10	0.52
1:B:710:MET:HE1	1:B:730[B]:ILE:HG22	1.92	0.51
1:B:40:GLU:HB2	1:B:43:LEU:HD12	1.91	0.51
1:B:176:ARG:CG	1:B:432:PHE:HB2	2.40	0.51
1:B:556:ASP:OD1	1:B:557:PHE:N	2.41	0.51
1:A:725:TYR:HB3	1:A:728:LEU:HB2	1.93	0.51
1:B:749:ASP:O	1:B:753:ARG:HG3	2.10	0.51
1:A:7:LEU:HD23	1:A:247:PRO:HG3	1.93	0.50
1:A:743:THR:OG1	1:A:746:GLN:HG3	2.11	0.50
1:A:679:GLU:HA	1:A:682:LYS:HD3	1.94	0.50
1:A:684:ASN:HB3	1:A:759:MET:HE1	1.93	0.50
1:A:710:MET:HE2	1:A:714[A]:MET:HE1	1.94	0.50
1:B:420:VAL:HG23	1:B:662:GLY:HA3	1.93	0.50
1:B:64:GLU:HB3	6:B:2084:HOH:O	2.11	0.50
1:B:550:GLU:HG2	6:B:2469:HOH:O	2.11	0.50
1:A:442:MET:HE1	1:A:536:LEU:HG	1.95	0.49
1:B:636:MET:HB2	1:B:639:ARG:HD2	1.94	0.49
1:A:189:ILE:HG21	1:A:234:LYS:HG3	1.93	0.49
1:A:460:PRO:HD3	1:A:485:TRP:CD1	2.47	0.49
1:A:33:TYR:CE1	1:A:35:PRO:HG3	2.47	0.49
1:B:704:HIS:CG	1:B:705:LEU:N	2.81	0.49
1:B:674:LEU:HB3	1:B:681:ARG:HG2	1.95	0.49
1:A:16:LYS:HG2	1:A:20[B]:GLN:OE1	2.13	0.49
1:A:230:LEU:HD23	1:A:233[C]:MET:CE	2.43	0.49
1:A:233[A]:MET:HE2	1:A:261:GLY:HA2	1.93	0.49
1:A:229:ALA:O	1:A:233[A]:MET:HG3	2.13	0.49
1:B:189:ILE:HG21	1:B:234:LYS:HG3	1.95	0.48
1:B:568:ASN:HB3	1:B:642[B]:LYS:CE	2.43	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:731:ARG:HD2	1:A:736:ALA:HB2	1.96	0.48
1:B:605:THR:HG23	1:B:632:GLY:HA2	1.95	0.48
1:B:81:ILE:HD11	1:B:509:SER:HB3	1.95	0.48
1:B:301:ALA:HA	1:B:304[A]:MET:HE2	1.96	0.48
1:B:365:PRO:HG3	1:B:394:THR:HG22	1.95	0.48
1:A:63:LEU:HD21	6:A:2117:HOH:O	2.13	0.48
1:A:250:ASN:ND2	1:A:250:ASN:C	2.67	0.48
1:B:111:PRO:HD2	6:B:2122:HOH:O	2.13	0.48
1:A:119:GLU:CD	1:A:134:LYS:HE2	2.39	0.48
1:B:116:LYS:HB2	1:B:117:MET:CE	2.43	0.48
1:A:134:LYS:O	1:A:138[B]:THR:HG23	2.14	0.47
1:B:569:ASN:O	1:B:571:PRO:HD3	2.14	0.47
1:B:230:LEU:HD23	1:B:233[C]:MET:CE	2.44	0.47
1:B:400:GLU:HG2	1:B:704:HIS:HB2	1.96	0.47
1:A:445:ALA:HA	1:A:482[B]:PHE:CE2	2.50	0.47
1:B:442:MET:HB2	1:B:479[A]:MET:CE	2.45	0.47
1:A:749:ASP:O	1:A:753:ARG:HG3	2.14	0.47
1:B:115:ILE:HG21	1:B:138[B]:THR:CG2	2.45	0.47
1:B:360:LEU:HD12	1:B:390:VAL:HG11	1.96	0.47
1:B:418:CYS:SG	2:B:1001:PYR:C	3.02	0.47
1:B:559:ILE:HD11	1:B:620:PRO:O	2.15	0.47
1:B:418:CYS:SG	2:B:1001:PYR:CB	3.03	0.47
1:B:547[A]:ILE:HB	1:B:556:ASP:HB3	1.97	0.47
1:A:409:PHE:HE2	1:A:422:PRO:HB2	1.80	0.47
1:A:730[A]:ILE:HD11	6:A:2573:HOH:O	2.15	0.47
1:B:324:ASP:OD2	4:B:9010:DTL:O3	2.30	0.47
1:B:696:GLU:CD	1:B:699:ILE:HD12	2.40	0.46
1:B:706:ASN:HD21	1:B:734:GLY:CA	2.28	0.46
1:A:123:LYS:HB3	1:A:123:LYS:HE2	1.39	0.46
1:A:389:LYS:HE3	1:A:393:ASP:OD2	2.14	0.46
1:B:708:ASN:OD1	1:B:730[A]:ILE:HG13	2.15	0.46
1:B:115:ILE:HG21	1:B:138[B]:THR:HG22	1.97	0.46
1:B:82:THR:O	1:B:239:LYS:HE3	2.16	0.46
1:B:306:ASP:OD1	1:B:359:THR:HG22	2.16	0.45
1:B:547[B]:ILE:HB	1:B:556:ASP:HB3	1.97	0.45
1:B:545:LYS:HB2	1:B:545:LYS:HZ3	1.80	0.45
1:B:645:VAL:HG23	1:B:759:MET:OXT	2.16	0.45
1:A:230:LEU:HA	1:A:233[C]:MET:HE3	1.98	0.45
1:A:445:ALA:HA	1:A:482[B]:PHE:CD2	2.51	0.45
1:B:63:LEU:O	1:B:63:LEU:HD12	2.16	0.45
1:B:95:LYS:O	1:B:96:ILE:HD13	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:165:LEU:HD21	1:A:490:TYR:HA	1.99	0.45
1:B:236:MET:HE1	1:B:260:PHE:O	2.16	0.45
1:B:423:MET:HB2	1:B:428[B]:GLN:HB3	1.99	0.45
1:A:81:ILE:HD11	1:A:509:SER:HB3	1.98	0.45
1:A:704:HIS:CG	1:A:705:LEU:N	2.84	0.45
1:B:438:LEU:HB2	1:B:526[B]:ILE:HG23	1.98	0.45
1:A:677:ASP:O	1:A:681:ARG:HG3	2.17	0.45
1:A:568:ASN:HB3	1:A:642[B]:LYS:CD	2.47	0.45
1:B:372:THR:HG23	1:B:398:GLN:HG3	1.99	0.44
1:A:606:ILE:HG22	1:A:607:THR:H	1.83	0.44
1:A:179:GLY:HA2	1:A:514:HIS:CE1	2.52	0.44
1:A:712:ARG:HG3	1:A:751[A]:ILE:CD1	2.48	0.44
1:B:250:ASN:HD21	1:B:253:GLU:H	1.62	0.44
1:B:328:SER:HB3	1:B:746:GLN:NE2	2.32	0.44
1:A:3:LEU:HA	1:A:7:LEU:HD12	2.00	0.44
1:A:255:ILE:HG12	1:A:283:LEU:HD13	2.00	0.44
1:A:406:ARG:CB	1:A:407:PRO:HD3	2.47	0.44
1:A:712:ARG:HG3	1:A:751[A]:ILE:HD11	1.99	0.44
1:B:677:ASP:CG	1:B:680:VAL:HG23	2.42	0.44
1:A:87:GLY:C	1:A:99:LEU:HD12	2.43	0.44
1:A:386:PHE:O	1:A:390:VAL:HG23	2.18	0.44
1:A:398:GLN:HB3	1:A:731:ARG:NH2	2.31	0.44
1:A:368:GLU:HA	1:A:369:PRO:C	2.43	0.43
1:A:62:LYS:O	1:A:66[B]:ARG:HG2	2.18	0.43
1:B:116:LYS:HB2	1:B:117:MET:HE1	1.99	0.43
1:A:38:GLY:C	1:A:382:ASN:HD22	2.26	0.43
1:A:356:PHE:O	1:A:371:MET:HE1	2.19	0.43
1:A:651:VAL:HG11	1:A:665:TYR:HB2	2.00	0.43
1:B:330:ASP:N	1:B:331:PRO:CD	2.81	0.43
1:B:628:PRO:HG3	6:B:2491:HOH:O	2.18	0.43
1:B:667:PHE:HB3	1:B:705:LEU:HD11	2.01	0.43
1:B:395:SER:OG	1:B:710:MET:HE3	2.18	0.43
1:B:562:GLU:HG2	6:B:2470:HOH:O	2.18	0.43
1:A:423:MET:HB2	1:A:428[B]:GLN:HB3	2.00	0.43
1:A:627:ALA:HA	1:A:628:PRO:HD3	1.84	0.43
1:A:423:MET:HA	1:A:428[A]:GLN:OE1	2.19	0.42
1:A:478:ARG:HA	1:A:478:ARG:NE	2.34	0.42
1:B:631:PRO:HD2	6:B:2505:HOH:O	2.18	0.42
1:A:668:SER:OG	1:A:708:ASN:HB2	2.19	0.42
1:B:117:MET:HG2	1:B:171:ALA:O	2.20	0.42
1:B:159:CYS:HB3	1:B:164:VAL:HG13	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:151:VAL:O	1:B:151:VAL:HG12	2.19	0.42
1:A:117:MET:N	1:A:117:MET:CE	2.71	0.42
1:B:454:LEU:O	1:B:455:LYS:C	2.63	0.42
1:B:743:THR:O	1:B:747:GLN:HG3	2.19	0.42
1:A:711:ASN:HB3	1:A:713[B]:GLU:OE1	2.20	0.42
1:B:716:LEU:HD21	1:B:751[A]:ILE:HG12	2.01	0.42
1:A:79:SER:HB3	1:A:110[A]:ILE:HD13	2.02	0.42
1:A:414:TYR:HA	1:A:424:ILE:HA	2.01	0.42
1:A:481[A]:HIS:O	1:A:484:ASP:HB2	2.20	0.42
1:A:680:VAL:HG12	1:A:684:ASN:ND2	2.35	0.42
1:B:330:ASP:N	1:B:331:PRO:HD3	2.35	0.41
1:A:453:LYS:HE3	1:A:614:LYS:O	2.20	0.41
1:A:606:ILE:HG22	1:A:607:THR:N	2.36	0.41
1:B:184:VAL:HG21	1:B:233[A]:MET:HE1	2.01	0.41
1:B:135:LYS:HE3	1:B:140:TYR:OH	2.20	0.41
1:B:233[A]:MET:CE	1:B:261:GLY:HA2	2.50	0.41
1:B:540:LYS:HA	6:B:2464:HOH:O	2.21	0.41
1:B:573:VAL:HG13	1:B:574:ASP:OD1	2.20	0.41
1:A:324:ASP:OD2	4:A:9010:DTL:O3	2.39	0.41
1:A:685:LEU:HG	1:A:689:MET:HE2	2.02	0.41
1:B:310:MET:O	1:B:314:MET:HG3	2.21	0.41
1:B:624:ARG:NH1	6:B:2496:HOH:O	2.53	0.41
1:A:215:GLN:OE1	1:A:215:GLN:N	2.49	0.41
1:A:645:VAL:HG13	5:A:9013:PG4:H41	2.01	0.41
1:A:656:PHE:CD1	1:A:703:GLN:HG3	2.56	0.41
1:B:545:LYS:HA	1:B:546:PRO:HD3	1.92	0.41
1:A:377[B]:GLU:OE2	1:A:694:HIS:ND1	2.52	0.41
1:B:250:ASN:HD22	1:B:253:GLU:H	1.63	0.41
1:A:476:MET:HE2	1:A:583:ARG:NE	2.36	0.40
1:B:696:GLU:OE2	1:B:699:ILE:HD12	2.22	0.40
1:B:119:GLU:OE1	1:B:134:LYS:HE3	2.20	0.40
1:B:211:VAL:O	1:B:212:ASN:C	2.64	0.40
1:B:712:ARG:NH1	1:B:716:LEU:HD11	2.36	0.40
1:A:328:SER:HB3	1:A:746:GLN:NE2	2.36	0.40
1:A:420:VAL:O	1:A:422:PRO:HD3	2.21	0.40
1:A:677:ASP:OD2	1:A:680:VAL:HG23	2.20	0.40
1:B:368:GLU:HA	1:B:369:PRO:C	2.47	0.40
1:B:409:PHE:HE2	1:B:422:PRO:HB2	1.86	0.40
1:B:717:ASP:OD2	1:B:721:ASN:ND2	2.55	0.40
1:A:442:MET:CE	1:A:479[A]:MET:HE2	2.52	0.40
1:B:162:SER:OG	1:B:164:VAL:HG12	2.21	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:712:ARG:HH11	1:B:712:ARG:HG2	1.85	0.40
1:A:645:VAL:HG23	1:A:759:MET:OXT	2.22	0.40
1:B:478:ARG:HA	1:B:478:ARG:NE	2.36	0.40
1:B:758[A]:SER:OG	1:B:759:MET:N	2.54	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	799/759 (105%)	763 (96%)	34 (4%)	2 (0%)	36	46
1	B	799/759 (105%)	759 (95%)	36 (4%)	4 (0%)	24	31
All	All	1598/1518 (105%)	1522 (95%)	70 (4%)	6 (0%)	30	38

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	733	SER
1	B	606	ILE
1	A	606	ILE
1	B	554	ALA
1	B	733	SER
1	B	571	PRO

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	680/638 (107%)	668 (98%)	12 (2%)	51	70
1	B	680/638 (107%)	671 (99%)	9 (1%)	61	77
All	All	1360/1276 (107%)	1339 (98%)	21 (2%)	57	75

All (21) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	117	MET
1	A	123	LYS
1	A	152	TYR
1	A	250	ASN
1	A	371	MET
1	A	435	ARG
1	A	456	MET
1	A	468[A]	ASP
1	A	468[B]	ASP
1	A	545	LYS
1	A	607	THR
1	A	733	SER
1	B	5	GLU
1	B	146[A]	GLN
1	B	146[B]	GLN
1	B	152	TYR
1	B	250	ASN
1	B	257	TRP
1	B	435	ARG
1	B	545	LYS
1	B	733	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	92	GLN
1	A	250	ASN
1	A	302	GLN
1	A	358	ASN
1	A	410	ASN
1	A	457	GLN
1	A	592	HIS

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Mol	Chain	Res	Type
1	A	706	ASN
1	A	727	GLN
1	B	92	GLN
1	B	250	ASN
1	B	269	GLN
1	B	302	GLN
1	B	358	ASN
1	B	410	ASN
1	B	457	GLN
1	B	589	GLN
1	B	601	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 18 ligands modelled in this entry, 7 are monoatomic - leaving 11 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	$\# Z > 2$
5	PG4	A	9013	-	12,12,12	0.67	0	11,11,11	0.43	0
4	DTL	B	9009	-	7,7,7	0.66	0	8,8,8	0.62	0
5	PG4	B	9013	-	12,12,12	0.71	0	11,11,11	0.43	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	DTL	A	9010	-	7,7,7	0.44	0	8,8,8	0.34	0
5	PG4	A	9012	-	12,12,12	0.61	0	11,11,11	0.55	0
4	DTL	A	9009	-	7,7,7	0.44	0	8,8,8	0.57	0
5	PG4	B	9011	-	12,12,12	0.60	0	11,11,11	0.46	0
4	DTL	B	9010	-	7,7,7	0.65	0	8,8,8	0.18	0
2	PYR	B	1001	-	5,5,5	1.13	0	3,6,6	1.98	1 (33%)
5	PG4	A	9011	-	12,12,12	0.52	0	11,11,11	0.53	0
2	PYR	A	1001	-	5,5,5	1.39	1 (20%)	3,6,6	2.01	1 (33%)

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
5	PG4	A	9013	-	-	3/10/10/10	-
4	DTL	B	9009	-	-	0/8/8/8	-
5	PG4	B	9013	-	-	8/10/10/10	-
4	DTL	A	9010	-	-	2/8/8/8	-
5	PG4	A	9012	-	-	7/10/10/10	-
4	DTL	A	9009	-	-	2/8/8/8	-
5	PG4	B	9011	-	-	6/10/10/10	-
4	DTL	B	9010	-	-	2/8/8/8	-
2	PYR	B	1001	-	-	0/4/4/4	-
5	PG4	A	9011	-	-	6/10/10/10	-
2	PYR	A	1001	-	-	1/4/4/4	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	PYR	OXT-C	-2.15	1.24	1.30

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	PYR	OXT-C-CA	3.29	122.70	113.59
2	B	1001	PYR	OXT-C-CA	3.18	122.40	113.59

There are no chirality outliers.

All (37) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	A	9010	DTL	C1-C2-C3-C4
4	B	9010	DTL	C1-C2-C3-C4
5	A	9011	PG4	O2-C3-C4-O3
5	A	9011	PG4	O3-C5-C6-O4
5	B	9011	PG4	O2-C3-C4-O3
5	A	9012	PG4	O2-C3-C4-O3
5	B	9013	PG4	O1-C1-C2-O2
5	B	9013	PG4	O4-C7-C8-O5
5	B	9011	PG4	O1-C1-C2-O2
4	A	9009	DTL	O3-C3-C4-O4
5	B	9011	PG4	O3-C5-C6-O4
4	A	9009	DTL	C2-C3-C4-O4
5	A	9012	PG4	O3-C5-C6-O4
5	B	9013	PG4	C4-C3-O2-C2
5	A	9011	PG4	C1-C2-O2-C3
5	A	9012	PG4	C3-C4-O3-C5
5	A	9013	PG4	O3-C5-C6-O4
5	B	9011	PG4	C3-C4-O3-C5
5	B	9013	PG4	C8-C7-O4-C6
5	B	9013	PG4	C5-C6-O4-C7
5	B	9013	PG4	C6-C5-O3-C4
5	A	9012	PG4	O1-C1-C2-O2
2	A	1001	PYR	OXT-C-CA-O3
5	B	9013	PG4	C1-C2-O2-C3
5	A	9013	PG4	C8-C7-O4-C6
5	A	9011	PG4	C3-C4-O3-C5
5	A	9012	PG4	C8-C7-O4-C6
5	A	9011	PG4	C4-C3-O2-C2
5	B	9011	PG4	C1-C2-O2-C3
5	B	9011	PG4	C5-C6-O4-C7
5	A	9012	PG4	C5-C6-O4-C7
5	A	9013	PG4	C5-C6-O4-C7
5	A	9011	PG4	C6-C5-O3-C4
4	A	9010	DTL	C1-C2-C3-O3
4	B	9010	DTL	C1-C2-C3-O3
5	B	9013	PG4	O3-C5-C6-O4
5	A	9012	PG4	C1-C2-O2-C3

There are no ring outliers.

6 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	A	9013	PG4	2	0
5	B	9013	PG4	1	0
4	A	9010	DTL	2	0
4	B	9010	DTL	2	0
2	B	1001	PYR	3	0
2	A	1001	PYR	2	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	759/759 (100%)	-0.46	4 (0%) 87 87	13, 47, 67, 115	39 (5%)
1	B	759/759 (100%)	-0.29	5 (0%) 84 85	16, 50, 81, 113	40 (5%)
All	All	1518/1518 (100%)	-0.37	9 (0%) 85 86	13, 49, 76, 115	79 (5%)

All (9) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	583	ARG	3.6
1	A	1	SER	3.1
1	B	3	LEU	3.0
1	A	123	LYS	2.9
1	B	481[A]	HIS	2.6
1	B	1	SER	2.4
1	A	723	GLU	2.3
1	A	481[A]	HIS	2.2
1	B	2	GLU	2.2

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 oligosaccharides 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
5	PG4	B	9013	13/13	0.65	0.15	104,105,106,106	0
3	NA	A	9004	1/1	0.66	0.24	95,95,95,95	0
5	PG4	A	9013	13/13	0.78	0.12	89,91,102,102	0
5	PG4	B	9011	13/13	0.78	0.12	87,89,92,93	0
3	NA	B	9003	1/1	0.78	0.20	77,77,77,77	0
4	DTL	B	9010	8/8	0.79	0.20	83,88,90,92	0
5	PG4	A	9011	13/13	0.81	0.12	88,91,95,95	0
3	NA	B	9006	1/1	0.82	0.12	77,77,77,77	0
4	DTL	A	9010	8/8	0.82	0.18	81,83,84,85	0
5	PG4	A	9012	13/13	0.82	0.13	74,76,93,95	0
4	DTL	B	9009	8/8	0.85	0.12	59,67,70,72	0
4	DTL	A	9009	8/8	0.88	0.10	71,75,77,77	0
3	NA	B	9001	1/1	0.93	0.06	70,70,70,70	0
2	PYR	B	1001	6/6	0.93	0.10	53,55,55,55	0
3	NA	A	9001	1/1	0.96	0.05	69,69,69,69	0
3	NA	A	9002	1/1	0.96	0.15	92,92,92,92	0
3	NA	B	9007	1/1	0.96	0.07	97,97,97,97	0
2	PYR	A	1001	6/6	0.97	0.05	40,42,44,45	0

6.5 Other polymers

There are no such residues in this entry.