



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:18 PM UTC

PDB ID : 4MJM / pdb_00004mjm
Title : Crystal Structure of the Inosine 5'-monophosphate Dehydrogenase, with a Short Internal Deletion of CBS Domain from Bacillus anthracis str. Ames
Authors : Kim, Y.; Makowska-Grzyska, M.; Gu, M.; Anderson, W.F.; Joachimiak, A.; Center for Structural Genomics of Infectious Diseases (CSGID)
Deposited on : 2013-09-03
Resolution : 2.25 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<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

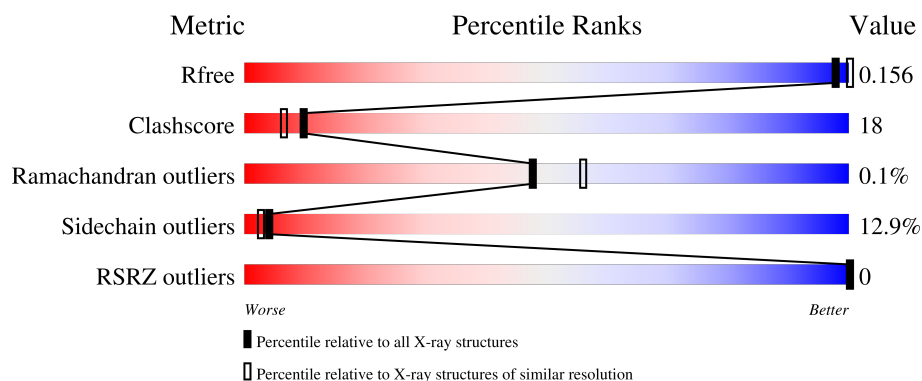
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.25 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	1898 (2.26-2.26)
Clashscore	190562	2005 (2.26-2.26)
Ramachandran outliers	187476	1965 (2.26-2.26)
Sidechain outliers	187428	1966 (2.26-2.26)
RSRZ outliers	180081	1898 (2.26-2.26)

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	385	
1	B	385	
1	C	385	
1	D	385	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 10136 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 Inosine-5'-monophosphate dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	328	Total	C	N	O	S	0	1	0
			2411	1514	421	463	13			
1	B	328	Total	C	N	O	S	0	0	0
			2408	1514	419	461	14			
1	C	329	Total	C	N	O	S	0	0	0
			2412	1517	419	462	14			
1	D	328	Total	C	N	O	S	0	0	0
			2409	1514	422	460	13			

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-2	SER	-	expression tag	UNP Q81W29
A	-1	ASN	-	expression tag	UNP Q81W29
A	0	ALA	-	expression tag	UNP Q81W29
A	95	GLY	-	linker	UNP Q81W29
B	-2	SER	-	expression tag	UNP Q81W29
B	-1	ASN	-	expression tag	UNP Q81W29
B	0	ALA	-	expression tag	UNP Q81W29
B	200	GLY	-	linker	UNP Q81W29
C	-2	SER	-	expression tag	UNP Q81W29
C	-1	ASN	-	expression tag	UNP Q81W29
C	0	ALA	-	expression tag	UNP Q81W29
C	95	GLY	-	linker	UNP Q81W29
D	-2	SER	-	expression tag	UNP Q81W29
D	-1	ASN	-	expression tag	UNP Q81W29
D	0	ALA	-	expression tag	UNP Q81W29
D	200	GLY	-	linker	UNP Q81W29

- Molecule 2 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	A	1	Total C O 4 2 2	0	0
2	B	1	Total C O 4 2 2	0	0
2	C	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0
2	D	1	Total C O 4 2 2	0	0

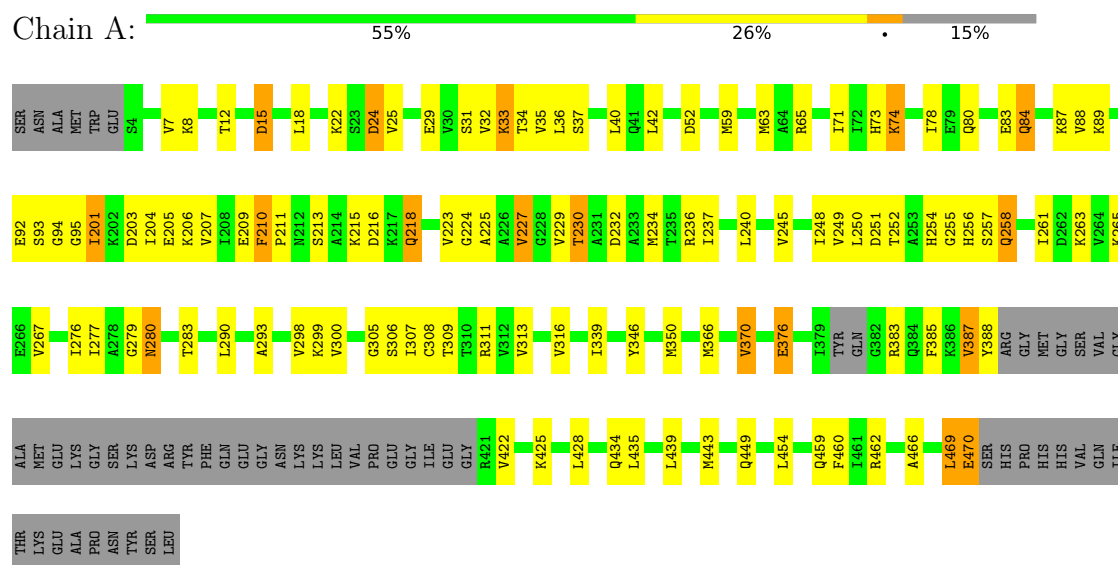
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	113	Total O 113 113	0	0
3	B	106	Total O 106 106	0	0
3	C	108	Total O 108 108	0	0
3	D	137	Total O 137 137	0	0

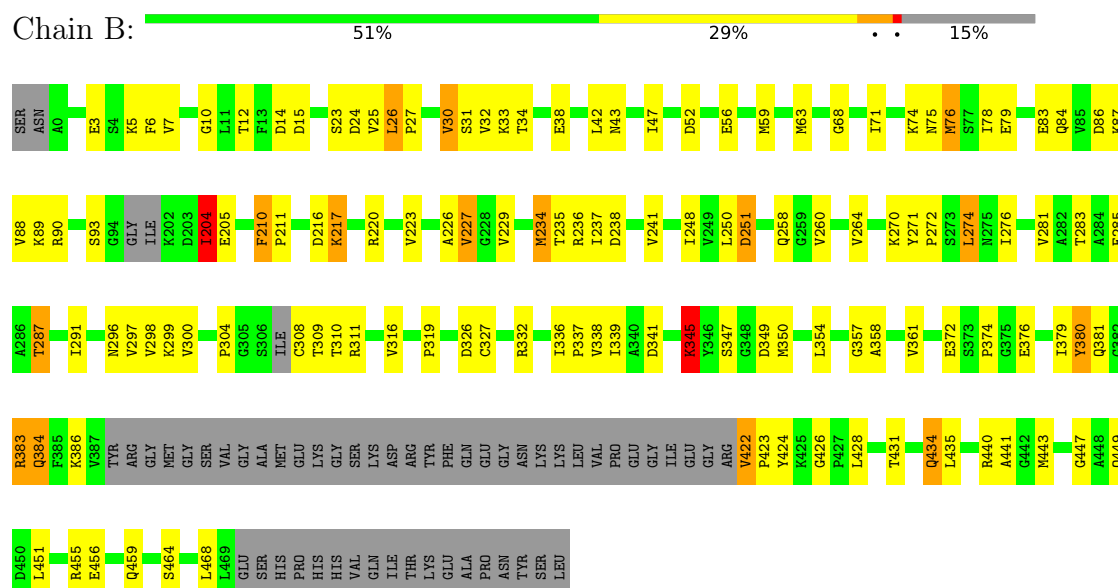
3 Residue-property plots [i](#)

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.

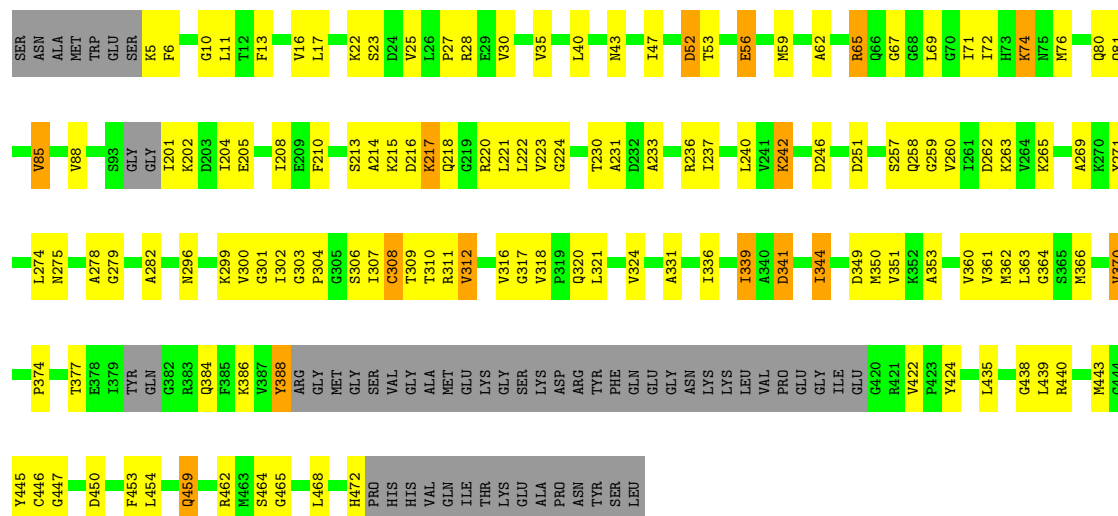
• Molecule 1: Inosine-5'-monophosphate dehydrogenase



• Molecule 1: Inosine-5'-monophosphate dehydrogenase



Chain C: 50% 29% 6% 15%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	84.33Å 84.25Å 84.31Å 110.01° 109.22° 109.19°	Depositor
Resolution (Å)	48.83 – 2.25 48.83 – 2.25	Depositor EDS
% Data completeness (in resolution range)	94.1 (48.83-2.25) 94.0 (48.83-2.25)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	6.29 (at 2.24Å)	Xtriage
Refinement program	PHENIX (phenix.refine: 1.8.2_1309)	Depositor
R, R_{free}	0.206 , 0.232 0.159 , 0.156	Depositor DCC
R_{free} test set	3951 reflections (4.66%)	wwPDB-VP
Wilson B-factor (Å ²)	40.4	Xtriage
Anisotropy	0.135	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.35 , 50.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.40$, $\langle L^2 \rangle = 0.22$	Xtriage

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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.

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Property	Value	Source
Estimated twinning fraction	0.077 for -k,-l,h+k+l 0.077 for h+k+l,-h,-k 0.457 for -l,-h,h+k+l 0.457 for -k,h+k+l,-h 0.078 for h+k+l,-l,-h 0.078 for -l,h+k+l,-k 0.078 for -h-k-l,k,h 0.078 for l,k,-h-k-l 0.078 for l,h,k 0.078 for k,l,h 0.078 for k,-h-k-l,l 0.078 for -h-k-l,h,l 0.078 for h,-h-k-l,k 0.078 for h,l,-h-k-l 0.377 for l,-h-k-l,h 0.077 for -h,h+k+l,-l 0.458 for -h-k-l,l,k 0.377 for -h,-l,-k 0.377 for k,h,-h-k-l 0.078 for -h,-k,h+k+l 0.077 for -l,-k,-h 0.377 for h+k+l,-k,-l 0.079 for -k,-h,-l	Xtriage
Reported twinning fraction	0.500 for -h-k-l,l,k	Depositor
Outliers	0 of 79280 reflections	Xtriage
F_o, F_c correlation	0.99	EDS
Total number of atoms	10136	wwPDB-VP
Average B, all atoms ($\text{\AA}^2$)	49.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.16% of the height of the origin peak. No significant pseudotranslation is detected.*

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: EDO

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	0.38	2/2438 (0.1%)	0.85	4/3290 (0.1%)
1	B	0.36	1/2436 (0.0%)	0.87	7/3287 (0.2%)
1	C	0.36	2/2440 (0.1%)	0.86	7/3294 (0.2%)
1	D	0.41	1/2436 (0.0%)	0.84	4/3286 (0.1%)
All	All	0.38	6/9750 (0.1%)	0.85	22/13157 (0.2%)

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	C	0	1

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	D	303	GLY	C-N	9.04	1.40	1.33
1	C	271	TYR	C-N	5.69	1.40	1.34
1	A	211	PRO	N-CD	5.38	1.55	1.47
1	C	272	PRO	N-CD	5.36	1.55	1.47
1	A	210	PHE	C-N	5.21	1.40	1.33
1	B	272	PRO	N-CD	5.20	1.55	1.47

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	233	ALA	N-CA-C	7.20	118.92	111.14
1	C	210	PHE	CA-C-N	7.11	126.64	119.24
1	C	210	PHE	C-N-CA	7.11	126.64	119.24

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	345	LYS	N-CA-C	6.65	118.32	111.14
1	A	387	VAL	N-CA-C	6.54	118.83	108.89
1	B	24	ASP	N-CA-C	-6.14	105.62	113.23
1	D	303	GLY	CA-C-N	-6.00	114.09	120.94
1	D	303	GLY	C-N-CA	-6.00	114.09	120.94
1	B	210	PHE	CA-C-N	5.80	125.90	119.87
1	B	210	PHE	C-N-CA	5.80	125.90	119.87
1	A	25	VAL	N-CA-C	5.68	116.03	107.80
1	D	306	SER	N-CA-C	-5.60	105.17	111.28
1	B	26	LEU	CA-C-N	5.55	126.78	119.84
1	B	26	LEU	C-N-CA	5.55	126.78	119.84
1	A	422	VAL	N-CA-C	5.55	113.94	107.89
1	B	204	ILE	N-CA-C	5.54	116.00	111.62
1	C	271	TYR	CA-C-N	-5.24	113.14	118.85
1	C	271	TYR	C-N-CA	-5.24	113.14	118.85
1	A	434	GLN	N-CA-C	-5.23	106.75	113.23
1	C	24	ASP	N-CA-C	-5.19	106.19	112.88
1	D	304	PRO	CA-N-CD	-5.06	104.92	112.00
1	C	270	LYS	N-CA-C	5.03	116.85	111.36

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	C	230	THR	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	2411	0	2482	92	0
1	B	2408	0	2475	97	0
1	C	2412	0	2487	88	0
1	D	2409	0	2482	89	0
2	A	12	0	18	0	0
2	B	4	0	6	0	0
2	C	4	0	6	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	12	0	18	1	0
3	A	113	0	0	7	0
3	B	106	0	0	18	0
3	C	108	0	0	13	0
3	D	137	0	0	9	0
All	All	10136	0	9974	345	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 18.

All (345) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:281:VAL:HG21	1:B:287:THR:HG23	1.48	0.94
1:B:455:ARG:HD3	3:B:704:HOH:O	1.77	0.83
1:D:76:MET:HB2	1:D:80:GLN:HG3	1.62	0.81
1:A:387:VAL:O	1:A:388:TYR:HB2	1.84	0.76
1:A:470:GLU:OE2	1:B:12:THR:HG21	1.86	0.75
1:D:47:ILE:HG12	1:D:69:LEU:HB3	1.69	0.74
1:D:65:ARG:HD2	1:D:210:PHE:CG	2.23	0.73
1:D:344:ILE:HG23	1:D:349:ASP:HB2	1.71	0.73
1:A:24:ASP:OD1	1:A:24:ASP:N	2.23	0.72
1:C:89:LYS:HE2	3:C:662:HOH:O	1.91	0.71
1:D:468:LEU:HB3	3:D:686:HOH:O	1.90	0.71
1:C:89:LYS:NZ	1:C:216:ASP:OD2	2.23	0.71
1:A:376:GLU:HG2	1:A:387:VAL:HG13	1.72	0.71
1:C:248:ILE:HG12	1:C:274:LEU:HD21	1.74	0.70
1:D:344:ILE:HD13	1:D:361:VAL:HG13	1.74	0.69
1:D:370:VAL:O	1:D:386:LYS:NZ	2.26	0.69
1:A:205:GLU:O	1:A:209:GLU:N	2.21	0.69
1:D:344:ILE:HD13	1:D:361:VAL:CG1	2.23	0.69
1:B:74:LYS:NZ	1:B:251:ASP:O	2.26	0.68
1:B:14:ASP:OD2	1:B:345:LYS:HE3	1.92	0.68
1:A:459:GLN:NE2	3:A:647:HOH:O	2.26	0.68
1:C:443:MET:HB2	1:C:448:ALA:HB3	1.76	0.68
1:D:43:ASN:HB2	1:D:67:GLY:HA3	1.77	0.67
1:D:230:THR:OG1	1:D:231:ALA:N	2.28	0.67
1:D:65:ARG:HD2	1:D:210:PHE:CD1	2.30	0.67
1:B:327:CYS:SG	3:B:647:HOH:O	2.51	0.67
1:A:230:THR:O	1:A:263:LYS:NZ	2.22	0.66
1:D:6:PHE:H	2:D:502:EDO:H21	1.60	0.66

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:366:MET:O	1:A:425:LYS:NZ	2.28	0.66
1:D:71:ILE:HD13	1:D:224:GLY:HA3	1.78	0.66
1:B:455:ARG:NH1	3:B:691:HOH:O	2.28	0.66
1:B:14:ASP:OD2	1:B:345:LYS:CE	2.44	0.66
1:C:304:PRO:HB2	1:C:342:GLY:HA3	1.76	0.66
1:D:216:ASP:OD2	1:D:220:ARG:HB2	1.96	0.65
1:A:18:LEU:HB2	1:B:316:VAL:HG12	1.79	0.65
1:D:472:HIS:NE2	3:D:697:HOH:O	2.30	0.65
1:A:258:GLN:NE2	3:A:648:HOH:O	2.29	0.65
1:D:72:ILE:HG13	1:D:85:VAL:HG22	1.79	0.65
1:D:301:GLY:HA3	1:D:341:ASP:OD2	1.97	0.65
1:C:5:LYS:HG2	1:C:6:PHE:N	2.12	0.65
1:B:379:ILE:HG12	1:B:384:GLN:HG2	1.77	0.64
1:B:379:ILE:HG12	1:B:384:GLN:CG	2.27	0.64
1:A:387:VAL:HG11	3:A:675:HOH:O	1.97	0.64
1:C:29:GLU:O	1:C:449:GLN:NE2	2.31	0.64
1:B:241:VAL:HG12	1:B:274:LEU:HD13	1.79	0.63
1:D:440:ARG:NH1	3:D:709:HOH:O	2.30	0.63
1:D:56:GLU:HB2	1:D:374:PRO:HG3	1.81	0.62
1:B:38:GLU:OE1	1:B:217:LYS:NZ	2.33	0.62
1:C:86:ASP:HA	1:C:89:LYS:HB2	1.82	0.62
1:C:80:GLN:NE2	3:C:619:HOH:O	2.33	0.62
1:D:282:ALA:HB3	1:D:302:ILE:HD12	1.82	0.62
1:D:311:ARG:HG2	1:D:317:GLY:HA3	1.82	0.62
1:A:74:LYS:NZ	1:A:251:ASP:OD1	2.32	0.62
1:B:229:VAL:HG12	1:B:229:VAL:O	1.99	0.62
1:B:234:MET:HE2	1:B:234:MET:H	1.65	0.62
1:B:350:MET:HE1	1:B:435:LEU:HB3	1.81	0.62
1:D:386:LYS:N	1:D:422:VAL:O	2.33	0.61
1:D:263:LYS:NZ	3:D:644:HOH:O	2.33	0.61
1:C:270:LYS:HG3	1:C:271:TYR:CD2	2.35	0.61
1:A:470:GLU:OE1	1:B:14:ASP:HB2	1.99	0.61
1:C:462:ARG:NH1	3:C:672:HOH:O	2.24	0.60
1:C:458:ALA:O	1:D:5:LYS:NZ	2.30	0.60
1:A:33:LYS:NZ	3:A:604:HOH:O	2.26	0.60
1:C:26:LEU:HD21	1:D:257:SER:HB2	1.82	0.60
1:A:305:GLY:O	1:D:472:HIS:ND1	2.35	0.60
1:C:310:THR:HB	1:C:314:ALA:HB3	1.84	0.60
1:D:308:CYS:SG	1:D:311:ARG:NH2	2.75	0.59
1:A:12:THR:OG1	1:A:15:ASP:OD1	2.20	0.59
1:B:300:VAL:N	1:B:339:ILE:O	2.30	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:89:LYS:HD2	1:C:220:ARG:HG2	1.84	0.59
1:A:34:THR:HG23	1:A:42:LEU:HB2	1.84	0.59
1:A:366:MET:HE1	1:A:435:LEU:HD21	1.84	0.59
1:B:300:VAL:O	1:B:341:ASP:N	2.28	0.59
1:A:29:GLU:O	1:A:449:GLN:NE2	2.35	0.59
1:D:282:ALA:CB	1:D:302:ILE:HD12	2.33	0.59
1:C:357:GLY:HA2	1:C:455:ARG:HD3	1.85	0.59
1:B:310:THR:N	3:B:667:HOH:O	2.29	0.58
1:D:13:PHE:CE2	1:D:344:ILE:HG13	2.39	0.58
1:D:275:ASN:HA	1:D:296:ASN:HD21	1.69	0.57
1:C:47:ILE:HG12	1:C:69:LEU:HB3	1.86	0.57
1:B:281:VAL:CG2	1:B:287:THR:HG23	2.27	0.57
1:A:94:GLY:N	1:A:203[B]:ASP:OD1	2.26	0.57
1:C:466:ALA:HB2	1:D:465:GLY:HA2	1.87	0.57
1:A:8:LYS:O	1:D:462:ARG:N	2.38	0.57
1:A:225:ALA:HB2	1:A:245:VAL:HG21	1.87	0.57
1:A:32:VAL:H	1:A:33:LYS:HE2	1.70	0.57
1:B:297:VAL:HG22	1:B:337:PRO:HG2	1.87	0.57
1:A:469:LEU:HD23	1:A:469:LEU:N	2.20	0.56
1:B:43:ASN:ND2	3:B:635:HOH:O	2.38	0.56
1:C:254:HIS:HB3	1:C:257:SER:HB3	1.87	0.56
1:A:92:GLU:HA	1:A:203[A]:ASP:OD2	2.05	0.56
1:A:261:ILE:HG23	1:A:293:ALA:HB2	1.87	0.56
1:B:33:LYS:HG2	1:B:43:ASN:HA	1.88	0.56
1:C:218:GLN:HB2	1:C:220:ARG:HE	1.71	0.56
1:D:339:ILE:HD11	1:D:362:MET:HB2	1.87	0.56
1:C:11:LEU:HD21	1:C:462:ARG:HD3	1.87	0.56
1:C:369:GLY:HA3	1:C:425:LYS:HG3	1.87	0.56
1:A:78:ILE:HG23	1:A:236:ARG:HA	1.87	0.56
1:B:10:GLY:HA3	1:B:319:PRO:HG2	1.88	0.56
1:C:310:THR:O	1:C:315:GLY:N	2.39	0.55
1:B:311:ARG:NH1	3:B:637:HOH:O	2.36	0.55
1:C:27:PRO:O	1:C:440:ARG:NH1	2.38	0.55
1:C:223:VAL:N	3:C:662:HOH:O	2.32	0.55
1:C:340:ALA:HB3	1:C:361:VAL:HG12	1.87	0.55
1:D:237:ILE:HA	1:D:240:LEU:HD12	1.89	0.55
1:C:94:GLY:HA3	1:C:220:ARG:HD3	1.87	0.55
1:C:362:MET:C	1:C:363:LEU:HD12	2.31	0.55
1:D:74:LYS:O	1:D:236:ARG:NH2	2.40	0.55
1:A:52:ASP:HA	1:A:73:HIS:CD2	2.42	0.55
1:A:279:GLY:HA3	1:A:299:LYS:HB2	1.89	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:460:PHE:H	1:B:5:LYS:HB2	1.71	0.55
1:B:386:LYS:HB3	1:B:422:VAL:HG13	1.89	0.55
1:A:313:VAL:HG12	1:D:438:GLY:HA3	1.89	0.55
1:A:227:VAL:HG21	1:A:237:ILE:HG12	1.87	0.54
1:B:291:ILE:HD13	1:B:336:ILE:HG13	1.89	0.54
1:B:347:SER:HB2	1:C:313:VAL:HA	1.89	0.54
1:D:331:ALA:O	1:D:336:ILE:N	2.37	0.54
1:C:223:VAL:HG12	3:C:662:HOH:O	2.07	0.54
1:C:232:ASP:OD1	1:C:232:ASP:N	2.33	0.54
1:D:27:PRO:HA	1:D:30:VAL:HG22	1.89	0.54
1:B:304:PRO:HD2	3:B:667:HOH:O	2.07	0.54
1:A:308:CYS:SG	1:A:309:THR:N	2.81	0.54
1:A:37:SER:HA	3:A:680:HOH:O	2.07	0.54
1:C:66:GLN:NE2	3:C:635:HOH:O	2.37	0.54
1:A:254:HIS:CG	1:A:257:SER:HB2	2.43	0.53
1:D:233:ALA:HB3	1:D:263:LYS:HE2	1.90	0.53
1:A:63:MET:HE2	1:A:428:LEU:HD21	1.89	0.53
1:A:462:ARG:NH2	1:B:6:PHE:O	2.40	0.53
1:B:345:LYS:HB2	1:B:349:ASP:OD2	2.08	0.53
1:D:301:GLY:HA2	1:D:320:GLN:HE22	1.73	0.53
1:C:227:VAL:HA	1:C:236:ARG:HH12	1.74	0.53
1:C:213:SER:N	3:C:674:HOH:O	2.42	0.53
1:A:466:ALA:HA	1:A:469:LEU:HG	1.90	0.53
1:B:52:ASP:OD1	1:B:52:ASP:N	2.42	0.52
1:C:11:LEU:HB3	1:C:321:LEU:HD23	1.91	0.52
1:B:204:ILE:HD12	1:B:205:GLU:H	1.74	0.52
1:A:93:SER:N	1:A:203[B]:ASP:OD1	2.42	0.52
1:D:279:GLY:HA3	1:D:299:LYS:HB3	1.91	0.52
1:A:88:VAL:HG11	1:A:223:VAL:HB	1.92	0.51
1:B:87:LYS:NZ	3:B:678:HOH:O	2.28	0.51
1:B:25:VAL:HB	1:B:447:GLY:HA2	1.92	0.51
1:A:223:VAL:HG13	1:A:245:VAL:HG23	1.91	0.51
1:B:380:TYR:C	1:B:380:TYR:CD2	2.89	0.51
1:C:344:ILE:HD13	1:C:344:ILE:H	1.75	0.51
1:C:27:PRO:HB3	1:C:440:ARG:HB3	1.92	0.51
1:A:201:ILE:HA	1:A:204:ILE:HG12	1.93	0.51
1:C:47:ILE:HG13	1:C:360:VAL:HG21	1.92	0.51
1:B:56:GLU:HG3	1:B:374:PRO:HG3	1.93	0.51
1:A:89:LYS:HD3	1:A:223:VAL:HG12	1.92	0.50
1:C:372:GLU:N	1:C:372:GLU:OE2	2.44	0.50
1:C:204:ILE:O	1:C:208:ILE:N	2.44	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:440:ARG:NH2	3:C:641:HOH:O	2.42	0.50
1:A:36:LEU:HB2	1:A:40:LEU:HB3	1.94	0.50
1:C:220:ARG:NH1	1:C:244:SER:OG	2.39	0.50
1:C:359:HIS:HA	1:C:455:ARG:HH11	1.75	0.50
1:C:375:GLY:HA2	1:C:387:VAL:HG12	1.94	0.50
1:B:298:VAL:HB	1:B:338:VAL:HG12	1.93	0.50
1:D:28:ARG:NH1	3:D:636:HOH:O	2.44	0.50
1:C:280:ASN:ND2	3:C:637:HOH:O	2.32	0.50
1:B:424:TYR:CZ	1:B:426:GLY:HA2	2.47	0.50
1:B:304:PRO:HG3	1:B:308:CYS:O	2.11	0.50
1:A:95:GLY:O	1:A:204:ILE:HD11	2.12	0.49
1:B:89:LYS:HD3	1:B:223:VAL:HG12	1.92	0.49
1:B:354:LEU:HB3	3:B:691:HOH:O	2.11	0.49
1:D:88:VAL:HG12	1:D:221:LEU:HB2	1.95	0.49
1:B:27:PRO:HG3	1:B:441:ALA:HA	1.94	0.49
1:C:76:MET:HE2	1:C:80:GLN:HB3	1.95	0.49
1:A:32:VAL:N	1:A:33:LYS:HE2	2.26	0.49
1:A:65:ARG:HD3	1:A:210:PHE:CE2	2.48	0.49
1:D:81:GLN:OE1	1:D:236:ARG:NH2	2.40	0.49
1:B:31:SER:N	1:B:449:GLN:O	2.33	0.49
1:D:443:MET:HG2	1:D:454:LEU:HD22	1.94	0.49
1:A:249:VAL:HG22	1:A:277:ILE:HB	1.95	0.49
1:C:454:LEU:O	1:C:458:ALA:HB2	2.13	0.49
1:D:271:TYR:HB3	1:D:274:LEU:HB2	1.95	0.49
1:A:305:GLY:H	1:A:311:ARG:CZ	2.26	0.48
1:B:86:ASP:HA	1:B:89:LYS:HB2	1.94	0.48
1:C:93:SER:HB3	1:C:220:ARG:HG3	1.94	0.48
1:C:348:GLY:HA3	1:D:312:VAL:HA	1.95	0.48
1:A:460:PHE:HD2	1:B:5:LYS:HD2	1.78	0.48
1:C:270:LYS:HA	3:C:708:HOH:O	2.13	0.48
1:A:22:LYS:NZ	3:A:636:HOH:O	2.43	0.48
1:B:379:ILE:HG12	1:B:384:GLN:HG3	1.96	0.48
1:B:451:LEU:HD22	3:B:691:HOH:O	2.12	0.48
1:D:76:MET:SD	1:D:81:GLN:HA	2.54	0.48
1:B:30:VAL:O	1:B:440:ARG:NH1	2.45	0.48
1:A:256:HIS:ND1	1:D:22:LYS:HA	2.29	0.47
1:A:470:GLU:CD	1:B:12:THR:HG21	2.39	0.47
1:A:15:ASP:OD1	1:A:15:ASP:N	2.48	0.47
1:B:251:ASP:OD2	1:B:251:ASP:N	2.47	0.47
1:A:387:VAL:HG22	1:A:388:TYR:N	2.28	0.47
1:A:305:GLY:HA3	1:A:311:ARG:HD2	1.95	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:203[B]:ASP:OD2	1:A:203[B]:ASP:N	2.45	0.47
1:C:251:ASP:OD2	1:C:280:ASN:ND2	2.48	0.47
1:B:361:VAL:HA	3:B:631:HOH:O	2.14	0.47
1:A:443:MET:HG2	1:A:454:LEU:HD13	1.96	0.47
1:A:234:MET:HE1	1:A:267:VAL:HA	1.96	0.47
1:B:84:GLN:O	1:B:88:VAL:HG23	2.15	0.47
1:D:446:CYS:SG	3:D:616:HOH:O	2.25	0.47
1:A:306:SER:HA	1:D:472:HIS:HA	1.96	0.46
1:B:42:LEU:HD13	1:B:68:GLY:HA2	1.96	0.46
1:B:358:ALA:O	1:B:455:ARG:NH1	2.49	0.46
1:A:248:ILE:HD11	1:A:276:ILE:HG12	1.98	0.46
1:C:37:SER:HB3	1:C:40:LEU:HB2	1.97	0.46
1:C:304:PRO:CG	1:C:307:ILE:HG22	2.45	0.46
1:D:344:ILE:HD13	1:D:361:VAL:HG11	1.96	0.46
1:A:255:GLY:HA2	1:A:280:ASN:O	2.15	0.46
1:D:25:VAL:HG22	1:D:447:GLY:HA2	1.97	0.46
1:A:84:GLN:HE21	1:A:84:GLN:N	2.13	0.46
1:B:63:MET:SD	1:B:428:LEU:HD21	2.56	0.46
1:B:71:ILE:HG22	3:B:617:HOH:O	2.16	0.46
1:D:265:LYS:HE2	1:D:269:ALA:HB2	1.98	0.46
1:A:387:VAL:CG2	1:A:388:TYR:N	2.78	0.46
1:C:54:VAL:HG13	1:C:367:PHE:HB2	1.98	0.46
1:C:378:GLU:HG3	1:C:383:ARG:O	2.15	0.46
1:A:240:LEU:HB3	1:A:245:VAL:HG11	1.97	0.46
1:A:280:ASN:C	1:A:280:ASN:HD22	2.24	0.46
1:A:71:ILE:HD13	1:A:224:GLY:HA3	1.97	0.46
1:B:210:PHE:HA	1:B:211:PRO:HD3	1.68	0.46
1:C:378:GLU:O	1:C:384:GLN:NE2	2.48	0.46
1:B:220:ARG:NH1	3:B:611:HOH:O	2.49	0.45
1:D:341:ASP:O	1:D:341:ASP:CG	2.59	0.45
1:C:46:LEU:CD2	1:C:363:LEU:HD13	2.46	0.45
1:C:250:LEU:HD12	1:C:264:VAL:HG22	1.98	0.45
1:C:284:ALA:HB1	1:C:330:GLU:HB2	1.99	0.45
1:D:214:ALA:HB1	1:D:222:LEU:HD12	1.99	0.45
1:B:78:ILE:HG12	1:B:236:ARG:HG3	1.98	0.45
1:C:359:HIS:HA	1:C:455:ARG:NH1	2.31	0.45
1:D:53:THR:HB	1:D:388:TYR:CD1	2.51	0.45
1:D:74:LYS:HG3	3:D:609:HOH:O	2.15	0.45
1:D:217:LYS:H	1:D:217:LYS:HG2	1.56	0.45
1:C:220:ARG:HH12	1:C:244:SER:CB	2.30	0.45
1:D:10:GLY:O	3:D:687:HOH:O	2.21	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:205:GLU:HA	1:D:208:ILE:HB	1.98	0.45
1:D:366:MET:CE	1:D:435:LEU:HD11	2.47	0.45
1:C:37:SER:HB3	1:C:40:LEU:H	1.81	0.45
1:A:84:GLN:O	1:A:88:VAL:HG23	2.17	0.44
1:C:255:GLY:HA3	1:C:280:ASN:H	1.82	0.44
1:D:213:SER:OG	1:D:215:LYS:HE3	2.17	0.44
1:B:14:ASP:OD2	1:B:345:LYS:HE2	2.18	0.44
1:C:347:SER:HB3	1:C:435:LEU:HD13	2.00	0.44
1:B:33:LYS:HB3	3:B:679:HOH:O	2.17	0.44
1:B:250:LEU:HD12	1:B:264:VAL:HG22	2.00	0.44
1:C:31:SER:N	1:C:449:GLN:O	2.36	0.44
1:D:13:PHE:HE2	1:D:344:ILE:HG13	1.81	0.44
1:D:351:VAL:HG22	1:D:439:LEU:HA	1.98	0.44
1:D:47:ILE:HG13	1:D:360:VAL:HG21	1.98	0.44
1:B:270:LYS:HE2	1:B:271:TYR:CD2	2.52	0.44
1:A:93:SER:OG	1:A:203[B]:ASP:OD2	2.30	0.44
1:D:13:PHE:CE1	1:D:353:ALA:HB2	2.53	0.44
1:B:380:TYR:CE2	1:B:381:GLN:HG2	2.53	0.44
1:B:443:MET:SD	3:B:616:HOH:O	2.61	0.44
1:D:52:ASP:OD2	1:D:52:ASP:N	2.49	0.44
1:D:384:GLN:HB2	1:D:424:TYR:HB2	2.00	0.43
1:A:213:SER:O	1:A:215:LYS:NZ	2.51	0.43
1:A:383:ARG:HB3	1:A:385:PHE:CE2	2.53	0.43
1:C:345:LYS:HG3	1:C:346:TYR:CD2	2.53	0.43
1:B:75:ASN:ND2	3:B:610:HOH:O	2.43	0.43
1:B:216:ASP:CG	1:B:217:LYS:H	2.26	0.43
1:D:459:GLN:HE21	1:D:459:GLN:HB3	1.57	0.43
1:A:59:MET:HE1	1:A:370:VAL:HG12	1.99	0.43
1:A:236:ARG:O	1:A:240:LEU:HG	2.18	0.43
1:B:283:THR:HG22	1:B:285:GLU:H	1.83	0.43
1:B:12:THR:HG23	1:B:14:ASP:H	1.84	0.43
1:B:86:ASP:O	1:B:90:ARG:HG2	2.19	0.43
1:D:300:VAL:HG11	1:D:324:VAL:HA	2.00	0.43
1:A:256:HIS:O	1:D:23:SER:HB3	2.19	0.43
1:B:12:THR:HG22	1:B:15:ASP:OD2	2.19	0.43
1:B:296:ASN:ND2	1:B:297:VAL:HG23	2.34	0.43
1:C:262:ASP:O	1:C:265:LYS:HG2	2.18	0.43
1:A:298:VAL:O	1:A:339:ILE:N	2.49	0.43
1:C:59:MET:HA	1:C:372:GLU:HB3	2.01	0.43
1:C:422:VAL:HA	1:C:423:PRO:HD3	1.89	0.43
1:D:307:ILE:O	1:D:307:ILE:HG13	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:366:MET:HE1	1:D:435:LEU:HD11	2.01	0.43
1:B:32:VAL:HA	3:B:616:HOH:O	2.19	0.42
1:B:34:THR:HG23	1:B:42:LEU:HB2	2.01	0.42
1:A:232:ASP:OD2	3:A:616:HOH:O	2.22	0.42
1:B:23:SER:OG	1:B:25:VAL:O	2.36	0.42
1:C:57:ALA:N	1:C:84:GLN:OE1	2.52	0.42
1:C:309:THR:N	3:C:649:HOH:O	2.52	0.42
1:B:299:LYS:HA	1:B:339:ILE:HB	2.00	0.42
1:B:383:ARG:H	1:B:383:ARG:HD2	1.84	0.42
1:B:422:VAL:HA	1:B:423:PRO:HD3	1.89	0.42
1:C:85:VAL:HG21	1:C:240:LEU:HD22	2.01	0.42
1:D:62:ALA:HA	1:D:65:ARG:HG3	2.01	0.42
1:B:248:ILE:HD11	1:B:276:ILE:HG23	2.00	0.42
1:C:19:VAL:HA	1:C:20:PRO:HD3	1.92	0.42
1:C:257:SER:O	1:C:261:ILE:HG13	2.19	0.42
1:D:278:ALA:O	1:D:299:LYS:N	2.37	0.42
1:A:218:GLN:HE21	1:A:218:GLN:HB3	1.65	0.42
1:B:383:ARG:H	1:B:383:ARG:CD	2.32	0.42
1:A:376:GLU:OE1	1:A:388:TYR:HB2	2.20	0.42
1:A:350:MET:HE2	1:A:439:LEU:HD22	2.01	0.42
1:B:270:LYS:CG	1:B:271:TYR:N	2.82	0.42
1:C:46:LEU:HD21	1:C:363:LEU:HD13	2.01	0.42
1:B:287:THR:HG21	1:B:327:CYS:HB3	2.01	0.42
1:C:464:SER:OG	3:C:663:HOH:O	2.22	0.42
1:B:380:TYR:CD2	1:B:380:TYR:O	2.73	0.41
1:C:288:LYS:NZ	1:C:334:HIS:HE1	2.18	0.41
1:D:259:GLY:HA2	1:D:262:ASP:HB2	2.00	0.41
1:C:350:MET:HE2	1:C:350:MET:HB3	1.95	0.41
1:D:223:VAL:HG23	1:D:246:ASP:H	1.85	0.41
1:A:311:ARG:HD3	1:D:17:LEU:HD21	2.01	0.41
1:B:76:MET:O	1:B:236:ARG:NH1	2.50	0.41
1:C:60:ALA:HB1	1:C:223:VAL:HG23	2.01	0.41
1:D:204:ILE:N	3:D:601:HOH:O	2.53	0.41
1:C:430:ASP:HA	1:C:433:HIS:CE1	2.54	0.41
1:D:450:ASP:CG	1:D:453:PHE:HB2	2.44	0.41
1:A:204:ILE:O	1:A:207:VAL:HB	2.21	0.41
1:B:79:GLU:O	1:B:83:GLU:N	2.52	0.41
1:C:448:ALA:HA	1:C:453:PHE:HD2	1.86	0.41
1:D:350:MET:HE2	1:D:350:MET:HB3	1.94	0.41
1:A:215:LYS:HE2	1:A:215:LYS:HB2	1.88	0.41
1:A:290:LEU:HD23	1:A:290:LEU:HA	1.91	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:316:VAL:HG21	1:D:445:TYR:HB2	2.01	0.41
1:A:206:LYS:HA	1:A:206:LYS:HD3	1.78	0.41
1:C:304:PRO:HG2	1:C:307:ILE:HG22	2.03	0.41
1:A:84:GLN:NE2	1:A:84:GLN:CA	2.84	0.41
1:A:261:ILE:HG22	1:A:265:LYS:HE3	2.03	0.41
1:B:59:MET:HA	1:B:372:GLU:HB3	2.01	0.41
1:B:434:GLN:HE21	1:B:434:GLN:HB3	1.75	0.41
1:D:344:ILE:HG23	1:D:349:ASP:CB	2.47	0.41
1:A:84:GLN:NE2	1:A:84:GLN:HA	2.36	0.41
1:B:74:LYS:HB3	1:B:226:ALA:O	2.21	0.41
1:C:257:SER:O	1:C:260:VAL:HG12	2.21	0.41
1:D:11:LEU:O	1:D:321:LEU:HB3	2.21	0.41
1:D:339:ILE:HB	1:D:360:VAL:HG13	2.03	0.41
1:B:227:VAL:HG21	1:B:237:ILE:HG12	2.03	0.41
1:B:332:ARG:HH12	1:C:2:TRP:HE1	1.67	0.41
1:C:27:PRO:O	1:C:30:VAL:HG22	2.21	0.41
1:C:64:ALA:HB1	1:C:221:LEU:HB2	2.03	0.41
1:A:92:GLU:HG3	1:A:203[A]:ASP:HB3	2.01	0.40
1:B:464:SER:O	1:B:468:LEU:HG	2.21	0.40
1:C:75:ASN:HA	3:C:704:HOH:O	2.20	0.40
1:C:89:LYS:NZ	1:C:220:ARG:HB3	2.35	0.40
1:D:450:ASP:OD2	1:D:453:PHE:N	2.43	0.40
1:A:300:VAL:N	1:A:339:ILE:O	2.52	0.40
1:A:263:LYS:O	1:A:267:VAL:HG23	2.22	0.40
1:B:237:ILE:HG23	1:B:248:ILE:HG21	2.04	0.40
1:B:299:LYS:NZ	3:B:676:HOH:O	2.54	0.40
1:D:242:LYS:NZ	1:D:242:LYS:HA	2.36	0.40
1:A:305:GLY:H	1:A:311:ARG:NE	2.18	0.40
1:B:357:GLY:HA2	1:B:455:ARG:HG2	2.04	0.40
1:D:22:LYS:O	1:D:447:GLY:HA3	2.21	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

The Analysed column shows the number of residues for which the backbone conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	323/385 (84%)	299 (93%)	24 (7%)	0	100	100
1	B	320/385 (83%)	292 (91%)	28 (9%)	0	100	100
1	C	323/385 (84%)	298 (92%)	25 (8%)	0	100	100
1	D	320/385 (83%)	296 (92%)	23 (7%)	1 (0%)	36	40
All	All	1286/1540 (84%)	1185 (92%)	100 (8%)	1 (0%)	48	56

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	364	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	253/299 (85%)	225 (89%)	28 (11%)	6	4
1	B	252/299 (84%)	222 (88%)	30 (12%)	5	3
1	C	253/299 (85%)	213 (84%)	40 (16%)	2	1
1	D	253/299 (85%)	221 (87%)	32 (13%)	4	3
All	All	1011/1196 (84%)	881 (87%)	130 (13%)	4	3

All (130) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	7	VAL
1	A	15	ASP
1	A	24	ASP
1	A	31	SER
1	A	33	LYS
1	A	35	VAL
1	A	74	LYS

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Mol	Chain	Res	Type
1	A	80	GLN
1	A	83	GLU
1	A	84	GLN
1	A	87	LYS
1	A	201	ILE
1	A	216	ASP
1	A	218	GLN
1	A	227	VAL
1	A	229	VAL
1	A	230	THR
1	A	250	LEU
1	A	252	THR
1	A	258	GLN
1	A	280	ASN
1	A	283	THR
1	A	307	ILE
1	A	346	TYR
1	A	370	VAL
1	A	376	GLU
1	A	469	LEU
1	A	470	GLU
1	B	3	GLU
1	B	7	VAL
1	B	26	LEU
1	B	30	VAL
1	B	47	ILE
1	B	76	MET
1	B	93	SER
1	B	204	ILE
1	B	217	LYS
1	B	227	VAL
1	B	234	MET
1	B	235	THR
1	B	238	ASP
1	B	251	ASP
1	B	258	GLN
1	B	260	VAL
1	B	274	LEU
1	B	287	THR
1	B	309	THR
1	B	326	ASP
1	B	345	LYS

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Mol	Chain	Res	Type
1	B	376	GLU
1	B	380	TYR
1	B	383	ARG
1	B	384	GLN
1	B	422	VAL
1	B	431	THR
1	B	434	GLN
1	B	456	GLU
1	B	459	GLN
1	C	4	SER
1	C	5	LYS
1	C	34	THR
1	C	53	THR
1	C	79	GLU
1	C	83	GLU
1	C	93	SER
1	C	201	ILE
1	C	205	GLU
1	C	206	LYS
1	C	208	ILE
1	C	212	ASN
1	C	215	LYS
1	C	220	ARG
1	C	223	VAL
1	C	230	THR
1	C	242	LYS
1	C	248	ILE
1	C	251	ASP
1	C	260	VAL
1	C	262	ASP
1	C	263	LYS
1	C	273	SER
1	C	306	SER
1	C	307	ILE
1	C	309	THR
1	C	313	VAL
1	C	316	VAL
1	C	344	ILE
1	C	345	LYS
1	C	362	MET
1	C	377	THR
1	C	378	GLU

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Mol	Chain	Res	Type
1	C	379	ILE
1	C	383	ARG
1	C	384	GLN
1	C	385	PHE
1	C	445	TYR
1	C	462	ARG
1	C	468	LEU
1	D	16	VAL
1	D	35	VAL
1	D	40	LEU
1	D	52	ASP
1	D	56	GLU
1	D	59	MET
1	D	65	ARG
1	D	74	LYS
1	D	85	VAL
1	D	201	ILE
1	D	202	LYS
1	D	217	LYS
1	D	218	GLN
1	D	242	LYS
1	D	251	ASP
1	D	258	GLN
1	D	260	VAL
1	D	308	CYS
1	D	309	THR
1	D	310	THR
1	D	312	VAL
1	D	316	VAL
1	D	318	VAL
1	D	339	ILE
1	D	341	ASP
1	D	344	ILE
1	D	363	LEU
1	D	370	VAL
1	D	377	THR
1	D	388	TYR
1	D	459	GLN
1	D	464	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (22) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	80	GLN
1	A	84	GLN
1	A	258	GLN
1	A	296	ASN
1	A	384	GLN
1	B	75	ASN
1	B	84	GLN
1	B	218	GLN
1	B	296	ASN
1	B	384	GLN
1	C	66	GLN
1	C	73	HIS
1	C	81	GLN
1	C	256	HIS
1	C	334	HIS
1	C	384	GLN
1	C	433	HIS
1	D	320	GLN
1	D	359	HIS
1	D	384	GLN
1	D	457	ASN
1	D	459	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](#)

8 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	EDO	B	501	-	3,3,3	0.42	0	2,2,2	0.38	0
2	EDO	D	502	-	3,3,3	0.33	0	2,2,2	0.32	0
2	EDO	A	502	-	3,3,3	0.32	0	2,2,2	0.49	0
2	EDO	A	503	-	3,3,3	0.42	0	2,2,2	0.12	0
2	EDO	A	501	-	3,3,3	0.42	0	2,2,2	0.98	0
2	EDO	D	501	-	3,3,3	0.36	0	2,2,2	0.78	0
2	EDO	D	503	-	3,3,3	0.29	0	2,2,2	0.28	0
2	EDO	C	501	-	3,3,3	0.25	0	2,2,2	0.62	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	EDO	B	501	-	-	1/1/1/1	-
2	EDO	D	502	-	-	0/1/1/1	-
2	EDO	A	502	-	-	0/1/1/1	-
2	EDO	A	503	-	-	1/1/1/1	-
2	EDO	A	501	-	-	0/1/1/1	-
2	EDO	D	501	-	-	1/1/1/1	-
2	EDO	D	503	-	-	0/1/1/1	-
2	EDO	C	501	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	503	EDO	O1-C1-C2-O2
2	D	501	EDO	O1-C1-C2-O2
2	C	501	EDO	O1-C1-C2-O2
2	B	501	EDO	O1-C1-C2-O2

There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	D	502	EDO	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	328/385 (85%)	-1.50	0 100 100	24, 41, 77, 128	1 (0%)
1	B	328/385 (85%)	-1.50	0 100 100	24, 43, 79, 143	0
1	C	329/385 (85%)	-1.44	0 100 100	29, 49, 91, 141	0
1	D	328/385 (85%)	-1.48	0 100 100	27, 45, 82, 161	0
All	All	1313/1540 (85%)	-1.48	0 100 100	24, 45, 84, 161	1 (0%)

There are no RSRZ outliers to report.

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
2	EDO	C	501	4/4	0.98	0.05	41,42,43,45	0
2	EDO	A	502	4/4	0.99	0.05	57,60,63,65	0
2	EDO	A	503	4/4	0.99	0.07	44,47,48,52	0
2	EDO	B	501	4/4	0.99	0.06	82,83,87,89	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
2	EDO	A	501	4/4	0.99	0.08	40,41,43,46	0
2	EDO	D	501	4/4	0.99	0.04	39,45,47,49	0
2	EDO	D	502	4/4	0.99	0.03	35,36,41,42	0
2	EDO	D	503	4/4	0.99	0.03	38,39,40,40	0

6.5 Other polymers [i](#)

There are no such residues in this entry.