



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 14, 2026 – 06:28 PM UTC

PDB ID : 1GK2 / pdb_00001gk2
Title : Histidine Ammonia-Lyase (HAL) Mutant F329G from *Pseudomonas putida*
Authors : Baedeker, M.; Schulz, G.E.
Deposited on : 2001-08-07
Resolution : 1.90 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

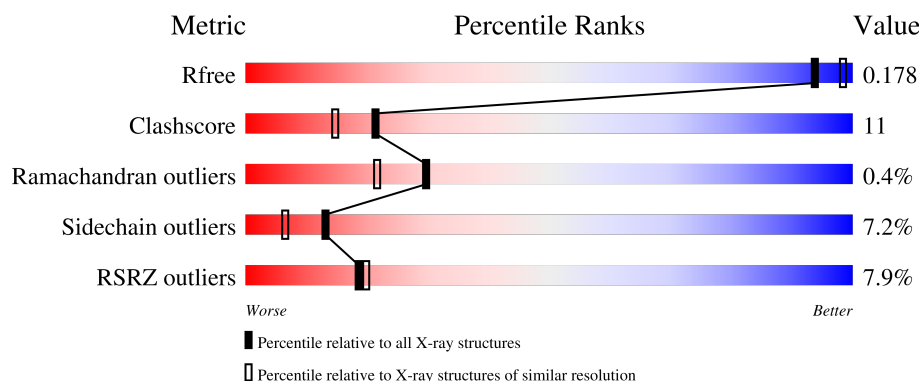
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.90 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	7789 (1.90-1.90)
Clashscore	190562	8410 (1.90-1.90)
Ramachandran outliers	187476	8333 (1.90-1.90)
Sidechain outliers	187428	8333 (1.90-1.90)
RSRZ outliers	180081	7790 (1.90-1.90)

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	509	8% 78% 19% .
1	B	509	8% 77% 20% .
1	C	509	8% 78% 20% ..
1	D	509	7% 77% 21% .

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard

residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	GOL	B	1510	-	-	X	-
2	GOL	C	1510	-	-	X	-
2	GOL	D	1510	-	-	X	-

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 16235 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 HISTIDINE AMMONIA-LYASE.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	509	Total	C	N	O	S	46	0	0
			3757	2353	670	718	16			
1	B	509	Total	C	N	O	S	39	0	0
			3757	2353	670	718	16			
1	C	509	Total	C	N	O	S	39	0	0
			3757	2353	670	718	16			
1	D	509	Total	C	N	O	S	39	0	0
			3757	2353	670	718	16			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	273	ALA	CYS	engineered mutation	UNP P21310
A	329	GLY	PHE	engineered mutation	UNP P21310
B	273	ALA	CYS	engineered mutation	UNP P21310
B	329	GLY	PHE	engineered mutation	UNP P21310
C	273	ALA	CYS	engineered mutation	UNP P21310
C	329	GLY	PHE	engineered mutation	UNP P21310
D	273	ALA	CYS	engineered mutation	UNP P21310
D	329	GLY	PHE	engineered mutation	UNP P21310

- Molecule 2 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



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		
2	C	1	Total	C	O	0	0
			6	3	3		
2	D	1	Total	C	O	0	0
			6	3	3		

- Molecule 3 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	C	1	Total O S 5 4 1	0	0
3	D	1	Total O S 5 4 1	0	0

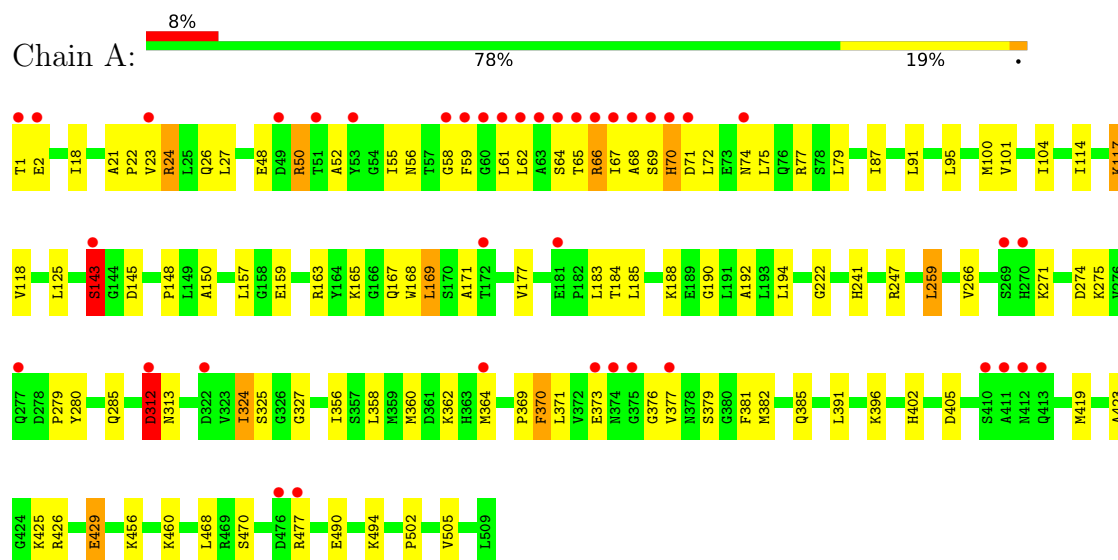
- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	289	Total O 289 289	0	0
4	B	298	Total O 298 298	0	0
4	C	286	Total O 286 286	0	0
4	D	290	Total O 290 290	0	0

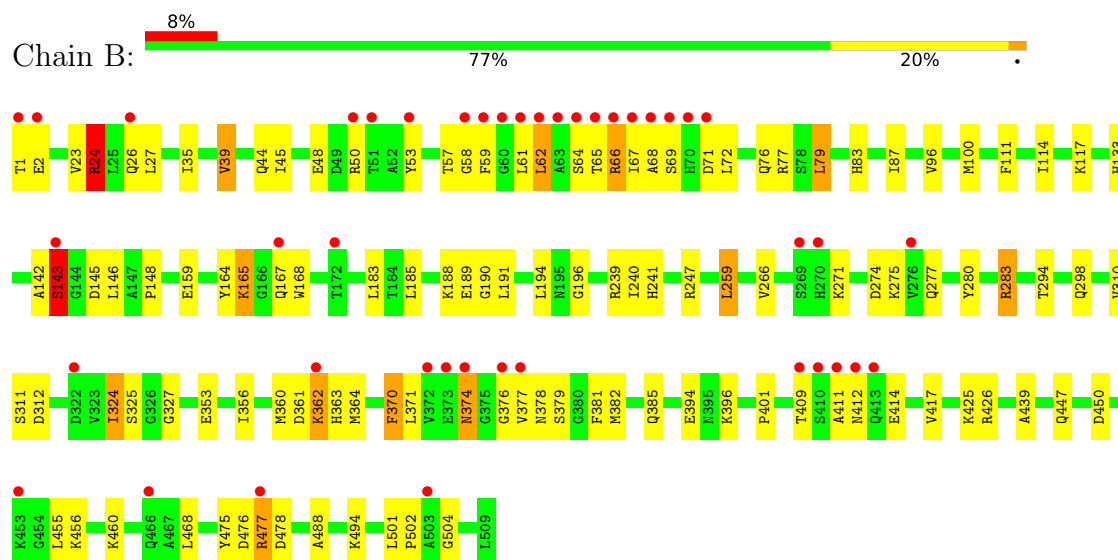
3 Residue-property plots

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.

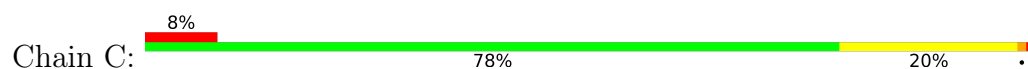
• Molecule 1: HISTIDINE AMMONIA-LYASE

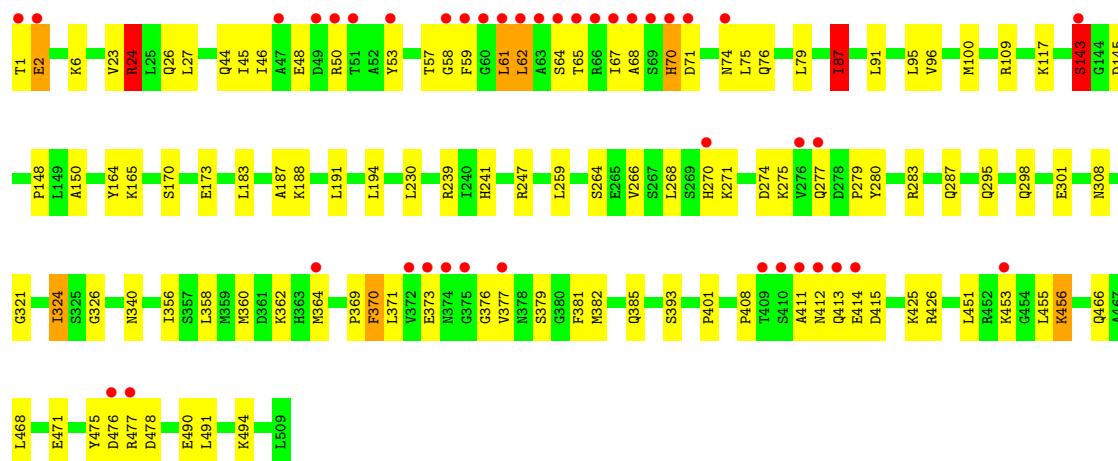


• Molecule 1: HISTIDINE AMMONIA-LYASE

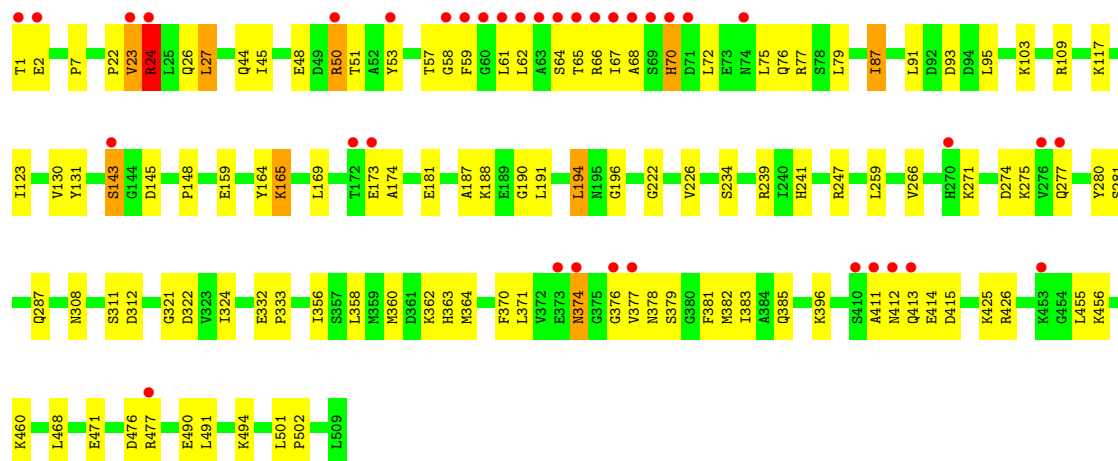
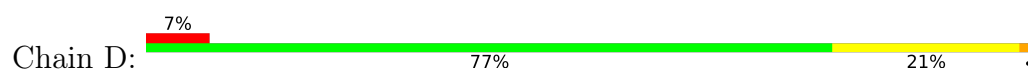


• Molecule 1: HISTIDINE AMMONIA-LYASE





• Molecule 1: HISTIDINE AMMONIA-LYASE



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	77.97Å 130.56Å 115.83Å 90.00° 91.00° 90.00°	Depositor
Resolution (Å)	25.00 – 1.90 25.00 – 1.90	Depositor EDS
% Data completeness (in resolution range)	77.0 (25.00-1.90) 76.9 (25.00-1.90)	Depositor EDS
R_{merge}	0.04	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.44 (at 1.91Å)	Xtriage
Refinement program	SHELX	Depositor
R, R_{free}	0.170 , 0.226 0.178 , 0.178	Depositor DCC
R_{free} test set	6936 reflections (4.99%)	wwPDB-VP
Wilson B-factor (Å ²)	12.1	Xtriage
Anisotropy	0.508	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.40 , 73.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Estimated twinning fraction	0.048 for h,-k,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	16235	wwPDB-VP
Average B, all atoms (Å ²)	19.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.49% 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: GOL, SO4

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	0.37	0/3814	1.09	8/5173 (0.2%)
1	B	0.37	0/3814	1.08	7/5173 (0.1%)
1	C	0.37	0/3814	1.07	9/5173 (0.2%)
1	D	0.37	0/3814	1.06	10/5173 (0.2%)
All	All	0.37	0/15256	1.07	34/20692 (0.2%)

There are no bond length outliers.

All (34) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	312	ASP	CA-CB-CG	8.37	120.97	112.60
1	A	241	HIS	CA-CB-CG	-7.43	106.37	113.80
1	D	241	HIS	CA-CB-CG	-7.25	106.55	113.80
1	D	143	SER	CA-C-N	7.17	132.52	121.12
1	D	143	SER	C-N-CA	7.17	132.52	121.12
1	B	241	HIS	CA-CB-CG	-7.16	106.64	113.80
1	A	143	SER	CA-C-N	6.76	130.62	121.83
1	A	143	SER	C-N-CA	6.76	130.62	121.83
1	C	87	ILE	CA-CB-CG1	6.26	121.05	110.40
1	C	143	SER	CA-C-N	6.15	129.47	121.23
1	C	143	SER	C-N-CA	6.15	129.47	121.23
1	B	24	ARG	CA-C-N	-5.98	113.20	122.59
1	B	24	ARG	C-N-CA	-5.98	113.20	122.59
1	C	241	HIS	CA-CB-CG	-5.98	107.82	113.80
1	A	370	PHE	CA-CB-CG	-5.91	107.89	113.80
1	C	370	PHE	CA-CB-CG	-5.87	107.93	113.80
1	A	24	ARG	CA-C-N	-5.83	113.02	122.81
1	A	24	ARG	C-N-CA	-5.83	113.02	122.81
1	B	312	ASP	CA-CB-CG	5.70	118.30	112.60
1	D	321	GLY	CA-C-N	5.70	130.52	122.09

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	D	321	GLY	C-N-CA	5.70	130.52	122.09
1	C	321	GLY	CA-C-N	5.59	131.62	122.07
1	C	321	GLY	C-N-CA	5.59	131.62	122.07
1	D	374	ASN	CA-C-N	5.52	127.08	120.13
1	D	374	ASN	C-N-CA	5.52	127.08	120.13
1	B	143	SER	CA-C-N	5.37	128.82	121.83
1	B	143	SER	C-N-CA	5.37	128.82	121.83
1	D	312	ASP	CA-CB-CG	5.14	117.74	112.60
1	C	24	ARG	CA-C-N	-5.11	114.23	122.82
1	C	24	ARG	C-N-CA	-5.11	114.23	122.82
1	A	21	ALA	N-CA-C	5.10	112.81	108.22
1	D	24	ARG	CA-C-N	-5.09	114.26	122.81
1	D	24	ARG	C-N-CA	-5.09	114.26	122.81
1	B	370	PHE	CA-CB-CG	-5.08	108.72	113.80

There are no chirality outliers.

There are no planarity outliers.

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	3757	0	3820	81	0
1	B	3757	0	3820	99	0
1	C	3757	0	3820	92	0
1	D	3757	0	3820	88	0
2	A	6	0	8	0	0
2	B	6	0	8	9	0
2	C	6	0	8	8	0
2	D	6	0	8	7	0
3	A	5	0	0	1	0
3	C	10	0	0	1	0
3	D	5	0	0	0	0
4	A	289	0	0	6	0
4	B	298	0	0	2	0
4	C	286	0	0	4	0
4	D	290	0	0	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	16235	0	15312	328	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 (328) 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:362:LYS:HD2	1:A:369:PRO:HG3	1.52	0.92
1:A:391:LEU:HD11	1:C:87:ILE:HD11	1.51	0.90
1:D:191:LEU:HG	2:D:1510:GOL:H12	1.57	0.87
1:A:377:VAL:HG13	1:C:75:LEU:HD13	1.59	0.85
1:B:376:GLY:HA3	1:D:62:LEU:HD22	1.62	0.80
1:A:117:LYS:HD3	1:A:183:LEU:HD21	1.64	0.79
1:C:279:PRO:HG2	1:C:358:LEU:HD11	1.64	0.79
1:C:360:MET:HE2	1:C:385:GLN:HB2	1.61	0.79
1:D:360:MET:HE2	1:D:385:GLN:HB2	1.65	0.79
1:A:360:MET:HE2	1:A:385:GLN:HB2	1.65	0.78
1:C:471:GLU:HG3	1:C:491:LEU:HD11	1.65	0.77
1:B:62:LEU:HD22	1:D:376:GLY:HA3	1.66	0.77
1:B:48:GLU:HB2	1:B:50:ARG:HE	1.50	0.77
1:A:59:PHE:O	1:C:376:GLY:HA2	1.86	0.76
1:A:66:ARG:O	1:A:67:ILE:HD13	1.87	0.74
1:D:53:TYR:HA	1:D:57:THR:OG1	1.86	0.74
1:B:239:ARG:NH2	1:C:324:ILE:HD11	2.02	0.73
1:D:45:ILE:HA	1:D:50:ARG:NE	2.03	0.72
1:D:490:GLU:HG3	1:D:494:LYS:HE3	1.71	0.72
1:C:379:SER:HB3	1:C:382:MET:HE2	1.69	0.72
1:D:44:GLN:O	1:D:48:GLU:HG3	1.90	0.72
1:B:69:SER:HA	1:B:72:LEU:HB2	1.73	0.71
1:B:266:VAL:HG23	1:B:456:LYS:O	1.91	0.71
1:B:377:VAL:HG13	1:D:75:LEU:HD13	1.70	0.71
1:D:1:THR:HG21	1:D:22:PRO:O	1.91	0.70
1:B:360:MET:HE2	1:B:385:GLN:HB2	1.74	0.68
1:A:279:PRO:HG2	1:A:358:LEU:HD11	1.75	0.68
1:A:266:VAL:HG23	1:A:456:LYS:O	1.94	0.68
1:B:188:LYS:H	2:B:1510:GOL:H32	1.59	0.68
1:B:259:LEU:HD22	1:B:460:LYS:HD3	1.77	0.67
1:A:68:ALA:HB1	1:A:70:HIS:NE2	2.09	0.67
1:C:170:SER:OG	1:C:173:GLU:HG3	1.94	0.67
1:C:53:TYR:HA	1:C:57:THR:OG1	1.94	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:266:VAL:HG22	1:C:455:LEU:HB3	1.77	0.67
1:D:364:MET:HA	1:D:364:MET:HE2	1.76	0.67
1:C:6:LYS:HB3	4:C:2012:HOH:O	1.93	0.66
1:C:266:VAL:HG23	1:C:456:LYS:O	1.95	0.66
1:B:143:SER:HB2	1:C:280:TYR:OH	1.96	0.66
1:A:376:GLY:HA3	1:C:62:LEU:HD22	1.78	0.66
1:B:379:SER:HB3	1:B:382:MET:HE2	1.79	0.65
1:B:188:LYS:HA	2:B:1510:GOL:H32	1.78	0.65
1:B:76:GLN:HE22	2:B:1510:GOL:H2	1.62	0.65
1:A:280:TYR:OH	1:D:143:SER:HB2	1.97	0.64
1:B:44:GLN:O	1:B:48:GLU:HG3	1.97	0.64
1:D:190:GLY:O	1:D:194:LEU:HG	1.98	0.64
1:A:67:ILE:HB	1:A:72:LEU:HD13	1.79	0.64
1:D:187:ALA:HA	2:D:1510:GOL:O3	1.98	0.64
1:A:59:PHE:CE2	1:A:75:LEU:HD11	2.33	0.64
1:B:409:THR:OG1	1:B:414:GLU:HB2	1.98	0.64
1:A:62:LEU:HA	1:A:65:THR:HB	1.78	0.63
1:D:259:LEU:HD22	1:D:460:LYS:HD3	1.80	0.63
1:B:114:ILE:HG13	1:B:189:GLU:HG2	1.81	0.63
2:C:1510:GOL:H32	4:C:2050:HOH:O	1.99	0.62
1:D:266:VAL:HG23	1:D:456:LYS:O	2.00	0.62
1:A:69:SER:HA	1:A:72:LEU:HB2	1.82	0.62
1:C:96:VAL:O	1:C:100:MET:HG3	2.00	0.61
1:A:379:SER:HB3	1:A:382:MET:HE2	1.81	0.61
1:D:2:GLU:HG2	1:D:24:ARG:HB3	1.82	0.61
1:A:490:GLU:HG3	1:A:494:LYS:HE3	1.82	0.60
1:C:59:PHE:CE2	1:C:75:LEU:HD11	2.36	0.60
1:B:277:GLN:HE21	1:B:283:ARG:CZ	2.15	0.60
1:C:490:GLU:HG3	1:C:494:LYS:HE3	1.83	0.60
1:C:364:MET:HE2	1:C:364:MET:HA	1.83	0.60
1:B:280:TYR:OH	1:C:143:SER:HB2	2.03	0.59
1:D:48:GLU:HB2	1:D:50:ARG:HG3	1.83	0.59
1:A:117:LYS:HG2	4:A:2087:HOH:O	2.01	0.59
1:B:142:ALA:O	1:B:143:SER:HB3	2.03	0.59
1:D:2:GLU:HA	1:D:24:ARG:O	2.03	0.58
1:D:27:LEU:HD13	1:D:123:ILE:HD11	1.85	0.58
1:C:143:SER:HB2	1:C:414:GLU:OE2	2.02	0.58
1:C:48:GLU:HB2	1:C:50:ARG:HE	1.67	0.58
1:D:266:VAL:HG22	1:D:455:LEU:HB3	1.85	0.58
1:D:59:PHE:CE2	1:D:75:LEU:HD11	2.37	0.58
1:C:62:LEU:HA	1:C:65:THR:HB	1.85	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:45:ILE:HA	1:D:50:ARG:HE	1.69	0.57
1:A:52:ALA:HB3	1:A:55:ILE:HB	1.86	0.57
1:B:360:MET:HE2	1:B:385:GLN:CB	2.33	0.57
1:D:362:LYS:HG3	1:D:363:HIS:N	2.18	0.57
1:D:266:VAL:HG22	1:D:455:LEU:CB	2.34	0.57
1:A:70:HIS:H	1:A:70:HIS:CD2	2.23	0.56
1:B:65:THR:HG22	1:B:66:ARG:O	2.04	0.56
1:C:76:GLN:OE1	2:C:1510:GOL:H11	2.05	0.56
1:C:70:HIS:H	1:C:70:HIS:CD2	2.24	0.56
2:D:1510:GOL:H32	4:D:2290:HOH:O	2.05	0.56
1:A:490:GLU:CG	1:A:494:LYS:HE3	2.36	0.56
1:D:76:GLN:HE22	2:D:1510:GOL:C3	2.19	0.56
1:C:68:ALA:HB1	1:C:70:HIS:NE2	2.21	0.55
1:B:240:ILE:HD11	1:C:324:ILE:CG2	2.35	0.55
1:B:188:LYS:N	2:B:1510:GOL:H32	2.22	0.55
1:D:188:LYS:HA	2:D:1510:GOL:H2	1.88	0.55
1:C:360:MET:HE2	1:C:385:GLN:CB	2.35	0.55
1:B:96:VAL:O	1:B:100:MET:HG3	2.06	0.55
1:A:59:PHE:HE2	1:A:75:LEU:HD11	1.71	0.55
1:A:356:ILE:O	1:A:360:MET:HG2	2.07	0.55
1:A:360:MET:HE2	1:A:385:GLN:CB	2.34	0.54
1:A:143:SER:HB2	1:D:280:TYR:OH	2.07	0.54
1:A:364:MET:HE2	1:A:364:MET:HA	1.90	0.54
1:B:377:VAL:CG1	1:D:75:LEU:HD13	2.37	0.54
1:C:356:ILE:O	1:C:360:MET:HG2	2.08	0.54
1:D:48:GLU:HB2	1:D:50:ARG:CG	2.38	0.54
1:B:356:ILE:O	1:B:360:MET:HG2	2.08	0.53
1:B:188:LYS:CA	2:B:1510:GOL:H32	2.38	0.53
1:B:376:GLY:CA	1:D:62:LEU:HD22	2.34	0.53
1:C:67:ILE:HG22	1:C:71:ASP:HB2	1.91	0.53
1:B:360:MET:HE3	1:B:370:PHE:HD1	1.72	0.53
1:C:164:TYR:CE2	1:C:165:LYS:HG3	2.43	0.53
1:C:266:VAL:HG22	1:C:455:LEU:CB	2.38	0.53
1:B:59:PHE:O	1:D:376:GLY:HA2	2.09	0.53
1:D:379:SER:HB3	1:D:382:MET:HE2	1.90	0.53
1:B:324:ILE:HD11	1:C:239:ARG:NH2	2.24	0.52
1:A:157:LEU:HD21	1:A:185:LEU:HG	1.91	0.52
1:B:362:LYS:HG3	1:B:363:HIS:N	2.24	0.52
1:B:143:SER:OG	1:B:146:LEU:HD11	2.09	0.52
1:C:76:GLN:HE22	2:C:1510:GOL:C3	2.22	0.52
1:B:79:LEU:O	1:B:83:HIS:HB2	2.10	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:145:ASP:HB3	1:D:148:PRO:HG2	1.91	0.51
1:D:396:LYS:NZ	4:D:2225:HOH:O	2.43	0.51
1:C:362:LYS:HD2	1:C:369:PRO:HG3	1.92	0.51
1:D:87:ILE:HD11	4:D:2045:HOH:O	2.10	0.51
1:D:411:ALA:O	1:D:412:ASN:HB2	2.11	0.51
1:B:48:GLU:CB	1:B:50:ARG:HE	2.22	0.51
1:D:70:HIS:H	1:D:70:HIS:CD2	2.27	0.51
1:B:298:GLN:CD	1:C:298:GLN:HB3	2.36	0.51
1:A:1:THR:O	1:A:1:THR:HG22	2.11	0.51
1:A:70:HIS:O	1:A:74:ASN:ND2	2.44	0.51
1:C:230:LEU:HD22	1:C:270:HIS:CD2	2.46	0.51
1:C:451:LEU:C	1:C:453:LYS:HD3	2.36	0.50
1:C:451:LEU:O	1:C:453:LYS:HD3	2.11	0.50
1:B:165:LYS:O	1:B:167:GLN:OE1	2.30	0.50
1:B:298:GLN:NE2	1:C:298:GLN:HB3	2.26	0.50
1:A:327:GLY:O	1:D:287:GLN:HG2	2.12	0.50
1:B:79:LEU:HD23	1:B:83:HIS:CE1	2.46	0.50
1:B:191:LEU:HG	2:B:1510:GOL:H11	1.93	0.50
1:B:240:ILE:HD13	1:C:326:GLY:HA3	1.94	0.50
1:B:401:PRO:HD3	1:C:401:PRO:HD3	1.93	0.50
1:C:360:MET:HE1	1:C:381:PHE:C	2.36	0.50
1:B:377:VAL:HG12	1:B:377:VAL:O	2.12	0.49
1:C:58:GLY:HA3	1:C:62:LEU:HB3	1.94	0.49
1:C:1:THR:O	1:C:1:THR:HG22	2.11	0.49
1:A:429:GLU:HG3	4:A:2237:HOH:O	2.11	0.49
1:B:44:GLN:HG2	1:B:50:ARG:HH21	1.78	0.49
1:D:471:GLU:HG3	1:D:491:LEU:HD11	1.94	0.49
1:B:277:GLN:HG3	1:B:283:ARG:NH2	2.27	0.49
1:C:277:GLN:HE21	1:C:283:ARG:CZ	2.26	0.49
1:B:240:ILE:HD11	1:C:324:ILE:HG21	1.94	0.49
1:A:177:VAL:HG23	4:A:2126:HOH:O	2.12	0.49
1:A:360:MET:HE1	1:A:381:PHE:C	2.37	0.49
1:B:45:ILE:HA	1:B:50:ARG:CZ	2.43	0.48
1:D:143:SER:HB2	1:D:414:GLU:OE2	2.13	0.48
1:B:360:MET:HE1	1:B:381:PHE:C	2.38	0.48
1:A:58:GLY:HA3	1:A:67:ILE:HD11	1.94	0.48
1:C:91:LEU:HD22	1:C:95:LEU:HD23	1.96	0.48
1:A:101:VAL:HA	1:A:104:ILE:HD12	1.96	0.48
1:B:48:GLU:HB2	1:B:50:ARG:HG3	1.95	0.48
1:B:277:GLN:HE21	1:B:283:ARG:NE	2.12	0.48
1:C:295:GLN:NE2	1:C:340:ASN:HB3	2.29	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:93:ASP:HB3	1:D:131:TYR:CD2	2.49	0.48
1:C:44:GLN:HG2	1:C:50:ARG:NH2	2.28	0.47
1:B:53:TYR:HA	1:B:57:THR:OG1	2.14	0.47
1:B:364:MET:HE2	1:B:364:MET:HA	1.95	0.47
1:C:475:TYR:OH	1:C:478:ASP:HA	2.15	0.47
1:B:62:LEU:HA	1:B:65:THR:HB	1.97	0.47
1:D:356:ILE:O	1:D:360:MET:HG2	2.14	0.47
1:A:74:ASN:HD22	1:A:74:ASN:H	1.62	0.47
1:C:183:LEU:HG	4:C:2108:HOH:O	2.15	0.47
1:A:159:GLU:O	1:A:171:ALA:HB2	2.15	0.47
1:C:2:GLU:HA	1:C:24:ARG:O	2.15	0.47
1:B:266:VAL:HG22	1:B:455:LEU:HB3	1.96	0.47
1:C:48:GLU:CD	1:C:50:ARG:HH21	2.22	0.47
1:D:68:ALA:HB1	1:D:70:HIS:CD2	2.50	0.47
1:A:58:GLY:HA3	1:A:62:LEU:HB3	1.97	0.47
1:A:266:VAL:HG23	1:A:456:LYS:C	2.39	0.47
1:D:1:THR:HG22	1:D:23:VAL:CG2	2.45	0.47
1:A:117:LYS:CD	1:A:183:LEU:HD21	2.40	0.47
1:A:324:ILE:HD11	1:D:239:ARG:NH2	2.30	0.47
1:A:402:HIS:HB2	1:A:423:ALA:HB2	1.96	0.47
1:C:408:PRO:HB3	1:C:415:ASP:HA	1.96	0.47
1:D:490:GLU:CG	1:D:494:LYS:HE3	2.44	0.47
1:A:114:ILE:HD11	1:A:118:VAL:HG11	1.96	0.46
1:A:169:LEU:N	1:A:169:LEU:HD23	2.30	0.46
1:A:188:LYS:O	1:A:188:LYS:HG3	2.15	0.46
1:D:58:GLY:HA3	1:D:62:LEU:HB3	1.98	0.46
1:D:24:ARG:HE	1:D:24:ARG:HA	1.80	0.46
1:A:145:ASP:HB3	1:A:148:PRO:HG2	1.97	0.46
1:B:294:THR:O	1:B:298:GLN:HG3	2.15	0.46
1:A:48:GLU:HB2	1:A:50:ARG:HG3	1.98	0.46
1:A:190:GLY:O	1:A:194:LEU:HG	2.16	0.46
1:A:18:ILE:HD13	1:A:101:VAL:HG11	1.96	0.46
1:A:377:VAL:HG12	1:A:377:VAL:O	2.16	0.46
1:B:1:THR:HG22	1:B:1:THR:O	2.16	0.46
1:B:142:ALA:HB3	1:B:417:VAL:HG12	1.97	0.46
1:B:411:ALA:O	1:B:412:ASN:HB2	2.15	0.46
1:C:150:ALA:HA	1:C:194:LEU:HD22	1.97	0.46
1:D:7:PRO:HB2	1:D:109:ARG:HG2	1.98	0.46
1:D:68:ALA:HB1	1:D:70:HIS:NE2	2.31	0.46
1:A:312:ASP:OD1	1:A:313:ASN:N	2.49	0.45
1:B:196:GLY:HA2	1:B:311:SER:O	2.16	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:504:GLY:HA2	4:B:2294:HOH:O	2.17	0.45
1:B:183:LEU:HG	4:B:2105:HOH:O	2.16	0.45
1:B:240:ILE:HD11	1:C:324:ILE:HG22	1.97	0.45
1:C:360:MET:HE3	1:C:370:PHE:HD1	1.82	0.45
1:D:45:ILE:HA	1:D:50:ARG:CZ	2.46	0.45
1:D:360:MET:HE1	1:D:381:PHE:C	2.41	0.45
1:A:68:ALA:O	1:A:71:ASP:N	2.50	0.45
1:C:46:ILE:HD11	1:C:187:ALA:CB	2.47	0.45
1:C:76:GLN:NE2	2:C:1510:GOL:H11	2.32	0.45
1:D:24:ARG:HE	1:D:24:ARG:CA	2.30	0.45
1:A:77:ARG:NH2	1:A:159:GLU:OE2	2.50	0.45
1:A:396:LYS:NZ	4:A:2217:HOH:O	2.50	0.45
1:B:327:GLY:O	1:C:287:GLN:HG2	2.17	0.45
1:A:100:MET:HE1	1:A:125:LEU:CB	2.47	0.45
1:B:361:ASP:OD2	1:B:364:MET:HG2	2.17	0.45
1:D:411:ALA:O	1:D:413:GLN:NE2	2.50	0.45
1:A:377:VAL:CG1	1:C:75:LEU:HD13	2.39	0.45
1:B:79:LEU:HD11	2:B:1510:GOL:H12	1.99	0.45
1:B:145:ASP:HB3	1:B:148:PRO:HG2	1.99	0.45
1:C:187:ALA:O	1:C:188:LYS:HB3	2.15	0.45
1:D:77:ARG:NH2	1:D:159:GLU:OE2	2.49	0.45
1:D:109:ARG:NH1	1:D:308:ASN:O	2.50	0.45
1:C:412:ASN:ND2	4:C:2228:HOH:O	2.50	0.45
1:D:62:LEU:O	1:D:65:THR:HB	2.16	0.45
1:D:360:MET:HE2	1:D:385:GLN:CB	2.41	0.45
1:A:58:GLY:CA	1:A:67:ILE:HD11	2.47	0.44
1:D:196:GLY:HA2	1:D:311:SER:O	2.16	0.44
1:D:266:VAL:HG23	1:D:456:LYS:C	2.42	0.44
1:A:56:ASN:O	1:A:66:ARG:HA	2.17	0.44
1:C:76:GLN:CD	2:C:1510:GOL:H11	2.43	0.44
1:C:377:VAL:HG12	1:C:377:VAL:O	2.18	0.44
1:D:44:GLN:C	1:D:50:ARG:HH21	2.26	0.44
1:D:374:ASN:HB2	1:D:378:ASN:ND2	2.31	0.44
1:B:447:GLN:O	1:B:450:ASP:HB2	2.16	0.44
1:C:283:ARG:NH1	3:C:1511:SO4:O1	2.50	0.44
1:B:76:GLN:NE2	2:B:1510:GOL:H2	2.31	0.44
1:B:77:ARG:NH2	1:B:159:GLU:OE2	2.50	0.44
1:C:79:LEU:HD21	1:C:191:LEU:CD2	2.48	0.44
1:C:145:ASP:HB3	1:C:148:PRO:HG2	2.00	0.44
1:A:313:ASN:ND2	3:A:1511:SO4:O3	2.50	0.44
1:B:24:ARG:HE	1:B:24:ARG:HA	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:58:GLY:HA3	1:B:62:LEU:HB3	2.00	0.44
1:D:169:LEU:HD22	1:D:173:GLU:CD	2.43	0.44
1:A:68:ALA:O	1:A:71:ASP:HB2	2.17	0.44
1:C:264:SER:O	1:C:268:LEU:HG	2.18	0.44
1:B:325:SER:OG	1:C:277:GLN:NE2	2.50	0.44
1:B:133:HIS:HB2	1:B:168:TRP:CZ3	2.52	0.43
1:A:74:ASN:HD22	1:A:74:ASN:N	2.16	0.43
1:B:360:MET:HE3	1:B:370:PHE:CD1	2.52	0.43
1:D:187:ALA:HA	2:D:1510:GOL:HO3	1.82	0.43
1:A:168:TRP:C	1:A:169:LEU:HD23	2.43	0.43
1:B:69:SER:O	1:B:72:LEU:HB3	2.18	0.43
1:C:411:ALA:O	1:C:413:GLN:NE2	2.50	0.43
1:C:412:ASN:HD21	1:D:412:ASN:HD21	1.67	0.43
1:B:35:ILE:O	1:B:39:VAL:HG23	2.19	0.43
1:B:188:LYS:H	2:B:1510:GOL:C3	2.27	0.43
1:B:239:ARG:HH21	1:C:324:ILE:HD11	1.80	0.43
1:A:58:GLY:CA	1:A:62:LEU:HB3	2.49	0.43
1:C:44:GLN:HG2	1:C:48:GLU:OE2	2.19	0.43
1:A:91:LEU:HD22	1:A:95:LEU:HD23	2.01	0.43
1:A:279:PRO:HG2	1:A:358:LEU:CD1	2.47	0.43
1:B:266:VAL:HG23	1:B:456:LYS:C	2.44	0.43
1:B:185:LEU:HD22	1:B:189:GLU:HB3	1.99	0.43
1:A:405:ASP:OD2	1:C:393:SER:OG	2.33	0.43
1:D:281:SER:OG	1:D:358:LEU:HD12	2.18	0.43
1:B:83:HIS:HA	1:D:383:ILE:HG13	2.01	0.42
1:B:111:PHE:HB2	1:B:310:VAL:HG11	2.00	0.42
1:B:439:ALA:HB1	1:B:488:ALA:HB3	2.01	0.42
1:C:70:HIS:O	1:C:74:ASN:ND2	2.51	0.42
1:C:279:PRO:CG	1:C:358:LEU:HD11	2.41	0.42
1:C:67:ILE:CG2	1:C:71:ASP:HB2	2.49	0.42
1:A:1:THR:HG21	1:A:22:PRO:O	2.18	0.42
1:B:69:SER:HA	1:B:72:LEU:CB	2.47	0.42
1:D:62:LEU:O	1:D:65:THR:N	2.52	0.42
1:A:163:ARG:HA	1:A:167:GLN:O	2.20	0.42
1:A:419:MET:HE3	1:C:393:SER:HB2	2.02	0.42
4:A:2217:HOH:O	1:B:396:LYS:NZ	2.51	0.42
1:B:68:ALA:O	1:B:71:ASP:HB2	2.18	0.42
1:B:277:GLN:HB3	1:B:283:ARG:HH21	1.83	0.42
1:C:188:LYS:O	1:C:188:LYS:HG3	2.20	0.42
1:C:191:LEU:HG	2:C:1510:GOL:O1	2.19	0.42
1:C:45:ILE:HA	1:C:50:ARG:CZ	2.49	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:222:GLY:HA3	1:A:285:GLN:OE1	2.20	0.42
1:B:164:TYR:CE2	1:B:165:LYS:HG3	2.54	0.42
1:B:48:GLU:HB2	1:B:50:ARG:NE	2.28	0.42
1:B:58:GLY:N	1:B:67:ILE:HG12	2.35	0.42
1:B:266:VAL:HG22	1:B:455:LEU:CB	2.50	0.42
1:B:475:TYR:OH	1:B:478:ASP:HA	2.19	0.42
1:C:44:GLN:HG2	1:C:50:ARG:HH21	1.85	0.42
1:C:283:ARG:O	1:C:283:ARG:HG3	2.20	0.42
1:A:163:ARG:HD2	4:A:2117:HOH:O	2.19	0.42
1:B:353:GLU:HA	1:B:356:ILE:HD12	2.02	0.42
1:B:363:HIS:O	1:B:364:MET:HE2	2.19	0.42
1:D:222:GLY:O	1:D:226:VAL:HG23	2.19	0.42
1:B:190:GLY:O	1:B:194:LEU:HG	2.19	0.41
1:B:374:ASN:HB2	1:B:378:ASN:ND2	2.35	0.41
1:D:59:PHE:HE2	1:D:75:LEU:HD11	1.82	0.41
1:C:266:VAL:HG23	1:C:456:LYS:C	2.45	0.41
1:D:66:ARG:O	1:D:67:ILE:HD13	2.19	0.41
1:D:332:GLU:N	1:D:333:PRO:HD2	2.36	0.41
1:A:188:LYS:HE2	1:A:192:ALA:HB2	2.01	0.41
1:A:150:ALA:HA	1:A:194:LEU:HD22	2.03	0.41
1:C:76:GLN:HE22	2:C:1510:GOL:H11	1.86	0.41
1:D:412:ASN:O	1:D:415:ASP:HB3	2.20	0.41
1:C:109:ARG:NH1	1:C:308:ASN:O	2.53	0.41
1:C:360:MET:HE3	1:C:370:PHE:CD1	2.56	0.41
1:D:259:LEU:CD2	1:D:460:LYS:HD3	2.50	0.41
1:A:325:SER:OG	1:D:277:GLN:NE2	2.53	0.41
1:D:62:LEU:CA	1:D:65:THR:HB	2.51	0.41
1:D:130:VAL:HG11	1:D:174:ALA:HB1	2.03	0.41
1:D:360:MET:HE3	1:D:370:PHE:HD1	1.85	0.41
1:C:61:LEU:HD23	1:C:61:LEU:N	2.36	0.41
1:C:76:GLN:HE22	2:C:1510:GOL:H31	1.85	0.41
1:A:502:PRO:O	1:A:505:VAL:HG23	2.21	0.40
1:D:1:THR:HG22	1:D:23:VAL:HG23	2.03	0.40
1:D:27:LEU:HD12	1:D:27:LEU:HA	1.94	0.40
1:B:501:LEU:HA	1:B:502:PRO:HD3	1.97	0.40
1:D:194:LEU:HD23	1:D:194:LEU:N	2.34	0.40
1:D:501:LEU:HG	1:D:502:PRO:HD2	2.04	0.40
1:A:75:LEU:O	1:A:75:LEU:HD12	2.21	0.40
1:B:476:ASP:HB2	1:B:477:ARG:H	1.68	0.40
1:C:45:ILE:HG12	1:C:50:ARG:NH1	2.36	0.40
1:D:51:THR:HG23	1:D:57:THR:HG22	2.04	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:187:ALA:O	1:D:188:LYS:HB3	2.21	0.40
1:D:76:GLN:NE2	2:D:1510:GOL:O3	2.50	0.40
1:D:164:TYR:CE2	1:D:165:LYS:HG3	2.56	0.40
1:A:66:ARG:C	1:A:67:ILE:HD13	2.47	0.40
1:A:100:MET:HE1	1:A:125:LEU:HB2	2.03	0.40
1:A:259:LEU:HD22	1:A:460:LYS:HD3	2.03	0.40
1:A:360:MET:HE3	1:A:370:PHE:HD1	1.86	0.40
1:D:91:LEU:HD22	1:D:95:LEU:HD23	2.03	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	507/509 (100%)	490 (97%)	14 (3%)	3 (1%)	21	13
1	B	507/509 (100%)	491 (97%)	13 (3%)	3 (1%)	21	13
1	C	507/509 (100%)	493 (97%)	12 (2%)	2 (0%)	30	22
1	D	507/509 (100%)	488 (96%)	18 (4%)	1 (0%)	43	36
All	All	2028/2036 (100%)	1962 (97%)	57 (3%)	9 (0%)	30	22

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	D	275	LYS
1	B	275	LYS
1	A	275	LYS
1	B	374	ASN
1	C	275	LYS
1	B	143	SER
1	A	66	ARG

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Mol	Chain	Res	Type
1	A	143	SER
1	C	143	SER

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	387/387 (100%)	358 (92%)	29 (8%)	12	6
1	B	387/387 (100%)	359 (93%)	28 (7%)	13	6
1	C	387/387 (100%)	361 (93%)	26 (7%)	15	7
1	D	387/387 (100%)	358 (92%)	29 (8%)	12	6
All	All	1548/1548 (100%)	1436 (93%)	112 (7%)	13	6

All (112) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	GLU
1	A	23	VAL
1	A	24	ARG
1	A	26	GLN
1	A	27	LEU
1	A	50	ARG
1	A	61	LEU
1	A	64	SER
1	A	70	HIS
1	A	79	LEU
1	A	87	ILE
1	A	117	LYS
1	A	165	LYS
1	A	169	LEU
1	A	184	THR
1	A	247	ARG
1	A	259	LEU
1	A	271	LYS
1	A	274	ASP

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Mol	Chain	Res	Type
1	A	312	ASP
1	A	324	ILE
1	A	371	LEU
1	A	373	GLU
1	A	425	LYS
1	A	426	ARG
1	A	429	GLU
1	A	468	LEU
1	A	470	SER
1	A	477	ARG
1	B	2	GLU
1	B	23	VAL
1	B	24	ARG
1	B	26	GLN
1	B	27	LEU
1	B	39	VAL
1	B	61	LEU
1	B	62	LEU
1	B	64	SER
1	B	66	ARG
1	B	79	LEU
1	B	87	ILE
1	B	117	LYS
1	B	165	LYS
1	B	247	ARG
1	B	259	LEU
1	B	271	LYS
1	B	274	ASP
1	B	283	ARG
1	B	324	ILE
1	B	362	LYS
1	B	371	LEU
1	B	394	GLU
1	B	425	LYS
1	B	426	ARG
1	B	468	LEU
1	B	477	ARG
1	B	494	LYS
1	C	2	GLU
1	C	23	VAL
1	C	24	ARG
1	C	26	GLN

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Mol	Chain	Res	Type
1	C	27	LEU
1	C	61	LEU
1	C	62	LEU
1	C	64	SER
1	C	70	HIS
1	C	87	ILE
1	C	117	LYS
1	C	247	ARG
1	C	259	LEU
1	C	271	LYS
1	C	274	ASP
1	C	301	GLU
1	C	324	ILE
1	C	371	LEU
1	C	373	GLU
1	C	425	LYS
1	C	426	ARG
1	C	456	LYS
1	C	466	GLN
1	C	468	LEU
1	C	476	ASP
1	C	477	ARG
1	D	23	VAL
1	D	24	ARG
1	D	26	GLN
1	D	27	LEU
1	D	50	ARG
1	D	61	LEU
1	D	64	SER
1	D	70	HIS
1	D	72	LEU
1	D	79	LEU
1	D	87	ILE
1	D	103	LYS
1	D	117	LYS
1	D	165	LYS
1	D	181	GLU
1	D	194	LEU
1	D	234	SER
1	D	247	ARG
1	D	271	LYS
1	D	274	ASP

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Mol	Chain	Res	Type
1	D	322	ASP
1	D	324	ILE
1	D	371	LEU
1	D	377	VAL
1	D	425	LYS
1	D	426	ARG
1	D	468	LEU
1	D	476	ASP
1	D	477	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (19) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	74	ASN
1	A	195	ASN
1	A	413	GLN
1	B	74	ASN
1	B	195	ASN
1	B	277	GLN
1	B	298	GLN
1	B	363	HIS
1	C	26	GLN
1	C	70	HIS
1	C	74	ASN
1	C	195	ASN
1	C	270	HIS
1	C	277	GLN
1	C	412	ASN
1	D	26	GLN
1	D	195	ASN
1	D	277	GLN
1	D	385	GLN

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 [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	GOL	B	1510	-	5,5,5	0.67	0	5,5,5	0.62	0
3	SO4	A	1511	-	4,4,4	0.72	0	6,6,6	0.08	0
3	SO4	C	1512	-	4,4,4	0.71	0	6,6,6	0.12	0
2	GOL	D	1510	-	5,5,5	0.65	0	5,5,5	0.45	0
2	GOL	C	1510	-	5,5,5	0.69	0	5,5,5	0.32	0
2	GOL	A	1510	-	5,5,5	0.65	0	5,5,5	0.40	0
3	SO4	D	1511	-	4,4,4	0.69	0	6,6,6	0.10	0
3	SO4	C	1511	-	4,4,4	0.72	0	6,6,6	0.08	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	GOL	A	1510	-	-	0/4/4/4	-
2	GOL	B	1510	-	-	4/4/4/4	-
2	GOL	D	1510	-	-	4/4/4/4	-
2	GOL	C	1510	-	-	2/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (10) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	B	1510	GOL	C1-C2-C3-O3
2	B	1510	GOL	O1-C1-C2-C3
2	C	1510	GOL	O1-C1-C2-C3
2	D	1510	GOL	O1-C1-C2-C3
2	D	1510	GOL	C1-C2-C3-O3
2	B	1510	GOL	O2-C2-C3-O3
2	C	1510	GOL	O1-C1-C2-O2
2	D	1510	GOL	O1-C1-C2-O2
2	D	1510	GOL	O2-C2-C3-O3
2	B	1510	GOL	O1-C1-C2-O2

There are no ring outliers.

5 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	1510	GOL	9	0
3	A	1511	SO4	1	0
2	D	1510	GOL	7	0
2	C	1510	GOL	8	0
3	C	1511	SO4	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 ⓘ

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	503/509 (98%)	0.07	40 (7%)	18	19	4, 13, 52, 89	0
1	B	504/509 (99%)	0.07	42 (8%)	17	18	4, 13, 51, 89	0
1	C	504/509 (99%)	0.07	41 (8%)	18	19	5, 13, 51, 89	0
1	D	504/509 (99%)	0.08	37 (7%)	21	22	4, 13, 52, 89	0
All	All	2015/2036 (98%)	0.07	160 (7%)	18	20	4, 13, 53, 89	0

All (160) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	1	THR	6.7
1	A	1	THR	6.6
1	C	63	ALA	6.6
1	C	70	HIS	6.4
1	D	61	LEU	6.2
1	A	70	HIS	6.1
1	C	276	VAL	6.1
1	D	70	HIS	6.1
1	A	59	PHE	6.0
1	B	1	THR	5.9
1	D	1	THR	5.8
1	B	69	SER	5.8
1	C	61	LEU	5.8
1	C	59	PHE	5.7
1	B	61	LEU	5.6
1	B	276	VAL	5.5
1	C	68	ALA	5.5
1	D	69	SER	5.5
1	B	143	SER	5.5
1	D	276	VAL	5.3
1	D	59	PHE	5.3

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Mol	Chain	Res	Type	RSRZ
1	A	60	GLY	5.2
1	C	53	TYR	5.2
1	B	68	ALA	5.1
1	B	59	PHE	5.1
1	D	63	ALA	5.0
1	A	63	ALA	5.0
1	B	60	GLY	5.0
1	D	67	ILE	5.0
1	A	411	ALA	4.9
1	B	411	ALA	4.9
1	A	61	LEU	4.8
1	A	69	SER	4.8
1	C	50	ARG	4.8
1	D	412	ASN	4.7
1	A	62	LEU	4.7
1	D	66	ARG	4.7
1	B	412	ASN	4.7
1	D	65	THR	4.6
1	C	270	HIS	4.6
1	D	50	ARG	4.5
1	B	63	ALA	4.5
1	A	68	ALA	4.3
1	D	410	SER	4.3
1	A	377	VAL	4.3
1	C	65	THR	4.2
1	B	270	HIS	4.2
1	D	411	ALA	4.2
1	A	270	HIS	4.1
1	B	64	SER	4.1
1	C	62	LEU	4.1
1	C	412	ASN	4.1
1	A	143	SER	4.0
1	C	375	GLY	4.0
1	A	67	ILE	4.0
1	B	67	ILE	3.9
1	B	66	ARG	3.8
1	C	66	ARG	3.8
1	D	477	ARG	3.8
1	A	65	THR	3.8
1	A	66	ARG	3.7
1	D	413	GLN	3.7
1	C	60	GLY	3.7

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Mol	Chain	Res	Type	RSRZ
1	D	53	TYR	3.7
1	B	477	ARG	3.7
1	C	67	ILE	3.7
1	B	53	TYR	3.7
1	C	69	SER	3.6
1	B	70	HIS	3.6
1	B	50	ARG	3.6
1	D	270	HIS	3.5
1	D	377	VAL	3.5
1	B	377	VAL	3.5
1	B	374	ASN	3.4
1	A	64	SER	3.4
1	A	412	ASN	3.4
1	D	2	GLU	3.4
1	A	410	SER	3.4
1	A	53	TYR	3.3
1	A	413	GLN	3.3
1	A	364	MET	3.2
1	D	64	SER	3.2
1	B	62	LEU	3.2
1	C	411	ALA	3.2
1	B	172	THR	3.2
1	C	377	VAL	3.2
1	D	374	ASN	3.2
1	D	58	GLY	3.1
1	C	71	ASP	3.1
1	B	65	THR	3.1
1	B	466	GLN	3.1
1	A	373	GLU	3.1
1	C	58	GLY	3.1
1	C	74	ASN	3.0
1	B	372	VAL	3.0
1	D	68	ALA	3.0
1	D	62	LEU	2.9
1	C	2	GLU	2.9
1	D	71	ASP	2.9
1	D	74	ASN	2.9
1	D	143	SER	2.8
1	B	167	GLN	2.8
1	B	373	GLU	2.8
1	C	64	SER	2.8
1	B	51	THR	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	71	ASP	2.8
1	C	477	ARG	2.7
1	C	143	SER	2.7
1	B	409	THR	2.7
1	D	373	GLU	2.7
1	C	453	LYS	2.7
1	D	60	GLY	2.6
1	B	413	GLN	2.6
1	C	51	THR	2.6
1	C	410	SER	2.6
1	A	374	ASN	2.6
1	C	374	ASN	2.6
1	A	312	ASP	2.6
1	A	269	SER	2.6
1	B	410	SER	2.6
1	A	477	ARG	2.6
1	D	376	GLY	2.5
1	A	23	VAL	2.5
1	D	277	GLN	2.5
1	A	322	ASP	2.5
1	A	476	ASP	2.5
1	A	74	ASN	2.5
1	B	2	GLU	2.5
1	D	23	VAL	2.5
1	C	364	MET	2.4
1	A	172	THR	2.4
1	B	362	LYS	2.4
1	B	58	GLY	2.4
1	A	2	GLU	2.4
1	C	373	GLU	2.4
1	C	49	ASP	2.4
1	C	409	THR	2.4
1	C	476	ASP	2.3
1	A	58	GLY	2.3
1	C	372	VAL	2.3
1	A	49	ASP	2.3
1	D	173	GLU	2.3
1	D	24	ARG	2.3
1	A	277	GLN	2.2
1	A	51	THR	2.2
1	B	71	ASP	2.2
1	B	503	ALA	2.1

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Mol	Chain	Res	Type	RSRZ
1	B	322	ASP	2.1
1	B	269	SER	2.1
1	B	453	LYS	2.1
1	D	453	LYS	2.1
1	B	26	GLN	2.1
1	A	181	GLU	2.1
1	A	375	GLY	2.1
1	C	413	GLN	2.1
1	B	376	GLY	2.0
1	C	47	ALA	2.0
1	C	414	GLU	2.0
1	C	277	GLN	2.0
1	D	172	THR	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 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($\text{\AA}^2$)	Q<0.9
3	SO4	A	1511	5/5	0.75	0.21	76,86,91,97	0
3	SO4	C	1512	5/5	0.80	0.23	73,103,104,117	0
3	SO4	C	1511	5/5	0.82	0.17	43,89,102,103	0
3	SO4	D	1511	5/5	0.84	0.14	40,81,83,103	0
2	GOL	B	1510	6/6	0.85	0.18	14,32,39,42	0
2	GOL	C	1510	6/6	0.86	0.20	16,46,51,72	0
2	GOL	D	1510	6/6	0.88	0.16	27,36,48,56	0
2	GOL	A	1510	6/6	0.91	0.11	25,34,38,43	0

6.5 Other polymers [i](#)

There are no such residues in this entry.