



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 7, 2026 – 01:22 AM UTC

PDB ID : 1P4D / pdb_00001p4d
Title : F factor TraI Relaxase Domain
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Deposited on : 2003-04-22
Resolution : 2.60 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

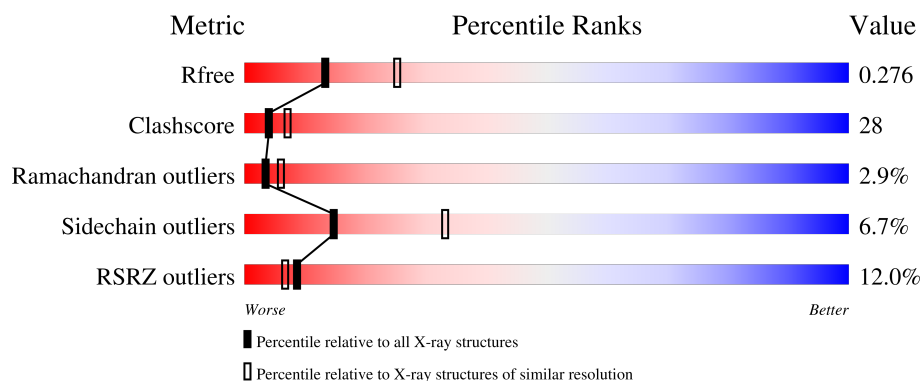
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.60 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	4008 (2.60-2.60)
Clashscore	190562	4347 (2.60-2.60)
Ramachandran outliers	187476	4277 (2.60-2.60)
Sidechain outliers	187428	4277 (2.60-2.60)
RSRZ outliers	180081	4008 (2.60-2.60)

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	330	2% 52% 25% 6% 17%
1	B	330	4% 54% 24% 5% 16%
1	C	330	25% 34% 43% 6% 16%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6656 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 TraI protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	275	Total	C	N	O	S	0	1	0
			2149	1331	387	418	13			
1	B	276	Total	C	N	O	S	0	0	0
			2148	1331	386	418	13			
1	C	276	Total	C	N	O	S	0	0	0
			2148	1331	386	418	13			

- Molecule 2 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	1	Total	Mg	0	0
			1	1		
2	B	1	Total	Mg	0	0
			1	1		
2	C	1	Total	Mg	0	0
			1	1		

- Molecule 3 is 1,2-ETHANEDIOL (CCD ID: EDO) (formula: C₂H₆O₂).



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

- Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	81	Total O 83 83	0	2

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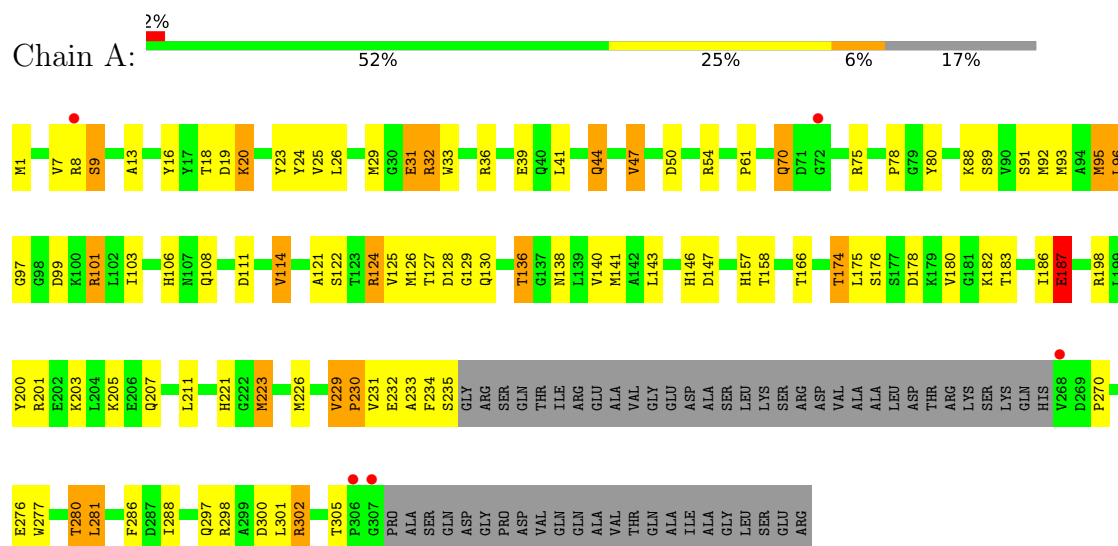
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	B	72	Total	O	0	0
			72	72		
4	C	9	Total	O	0	0
			9	9		

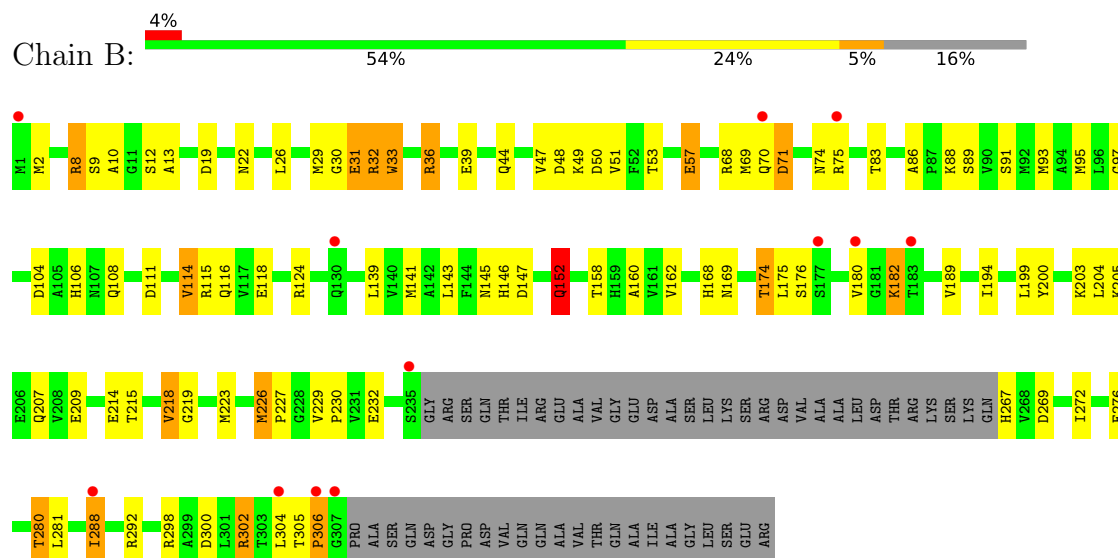
3 Residue-property plots [i](#)

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: TraI protein

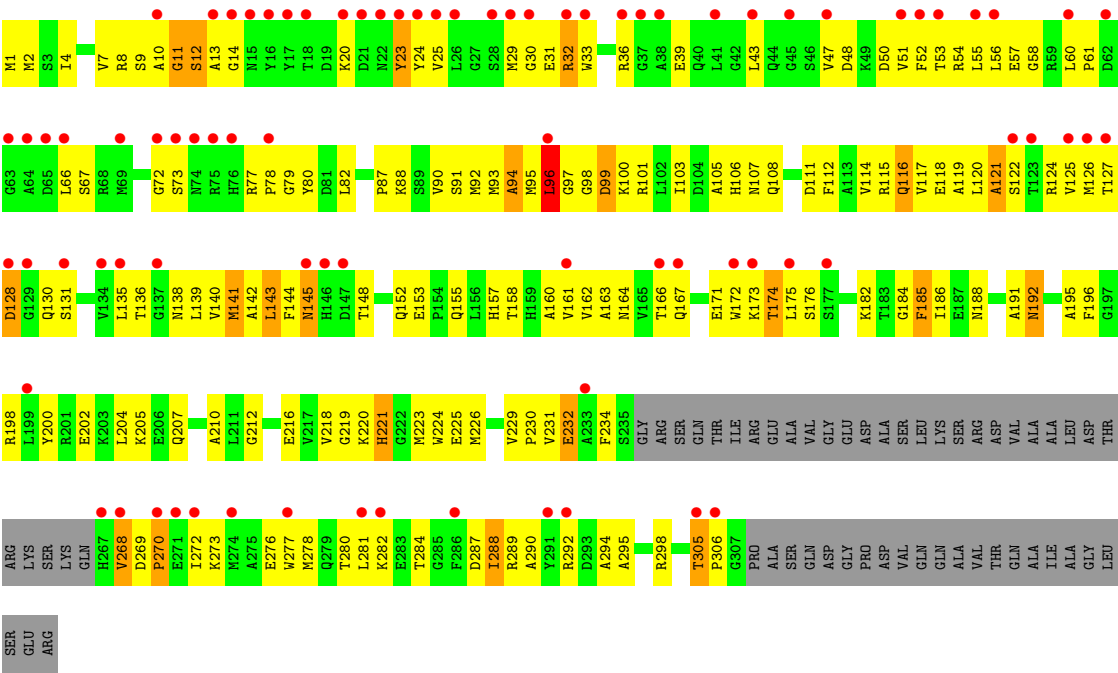


• Molecule 1: TraI protein



• Molecule 1: TraI protein





4 Data and refinement statistics

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, α , β , γ	128.22Å 128.22Å 121.18Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	25.00 – 2.60 25.00 – 2.60	Depositor EDS
% Data completeness (in resolution range)	96.9 (25.00-2.60) 96.7 (25.00-2.60)	Depositor EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	5.44 (at 2.50Å)	Xtriage
Refinement program	CNS 1.1	Depositor
R, R_{free}	0.233 , 0.278 0.233 , 0.276	Depositor DCC
R_{free} test set	3441 reflections (8.56%)	wwPDB-VP
Wilson B-factor (Å ²)	52.1	Xtriage
Anisotropy	0.085	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 67.9	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.023 for -h,-k,l	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	6656	wwPDB-VP
Average B, all atoms (Å ²)	67.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.89% 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: MG, 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	0.52	0/2187	0.99	10/2950 (0.3%)
1	B	0.48	0/2187	0.95	8/2951 (0.3%)
1	C	0.35	0/2187	0.81	3/2951 (0.1%)
All	All	0.46	0/6561	0.92	21/8852 (0.2%)

There are no bond length outliers.

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
1	B	33	TRP	N-CA-C	-7.31	101.10	110.53
1	A	229	VAL	CA-C-N	7.08	128.69	119.84
1	A	229	VAL	C-N-CA	7.08	128.69	119.84
1	A	18	THR	N-CA-C	6.66	123.38	114.12
1	B	189	VAL	N-CA-C	-6.24	104.67	110.53
1	A	96	LEU	N-CA-C	-6.04	106.55	114.04
1	B	226	MET	N-CA-C	-5.98	102.51	109.93
1	A	187	GLU	N-CA-C	-5.96	104.86	111.36
1	B	57	GLU	N-CA-C	-5.86	105.96	113.23
1	C	200	TYR	N-CA-C	-5.54	104.10	111.24
1	A	47	VAL	N-CA-C	5.50	116.03	107.99
1	C	116	GLN	N-CA-C	-5.41	104.93	112.45
1	A	103	ILE	N-CA-C	-5.41	105.23	110.42
1	B	86	ALA	CA-C-N	-5.40	114.68	120.03
1	B	86	ALA	C-N-CA	-5.40	114.68	120.03
1	A	95	MET	N-CA-C	5.22	119.75	113.17
1	C	141	MET	N-CA-C	5.21	118.24	109.95
1	B	175	LEU	N-CA-C	-5.21	105.21	112.45
1	A	175	LEU	N-CA-C	-5.17	105.33	111.69
1	B	152	GLN	N-CA-C	5.11	118.49	111.54
1	A	140	VAL	N-CA-C	-5.06	100.37	107.75

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2149	0	2086	91	0
1	B	2148	0	2081	95	0
1	C	2148	0	2081	173	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
3	A	20	0	30	3	0
3	B	20	0	30	0	0
3	C	4	0	6	0	0
4	A	83	0	0	4	0
4	B	72	0	0	0	0
4	C	9	0	0	2	0
All	All	6656	0	6314	356	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 28.

All (356) 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:29:MET:HE3	1:A:31:GLU:HG2	1.19	1.12
1:C:29:MET:HE3	1:C:31:GLU:HG2	1.25	1.10
1:A:88:LYS:HG3	1:A:92:MET:HE2	1.43	1.00
1:B:8:ARG:HH11	1:B:8:ARG:H	1.12	0.98
1:A:136:THR:HG21	1:A:166:THR:OG1	1.63	0.98
1:B:305:THR:HB	1:B:306:PRO:HD2	1.45	0.95
1:C:92:MET:HE1	1:C:231:VAL:HA	1.50	0.93
1:A:198:ARG:HD2	1:A:201:ARG:HH11	1.39	0.87
1:C:33:TRP:HE1	1:C:47:VAL:HA	1.41	0.86
1:C:268:VAL:HG12	1:C:269:ASP:H	1.40	0.85
1:B:218:VAL:HG12	1:B:223:MET:HE3	1.58	0.83

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:269:ASP:HB3	1:B:272:ILE:HG12	1.59	0.83
1:B:93:MET:HE1	1:B:288:ILE:HG12	1.58	0.83
1:B:93:MET:HB2	1:B:226:MET:HE1	1.61	0.81
1:B:8:ARG:HG2	1:B:9:SER:H	1.46	0.79
1:A:101:ARG:NH1	1:A:211:LEU:HD22	1.98	0.79
1:B:8:ARG:HG2	1:B:9:SER:N	1.98	0.79
1:B:219:GLY:HA3	1:B:223:MET:HE2	1.63	0.79
1:A:111:ASP:HA	3:A:612:EDO:H21	1.64	0.78
1:C:93:MET:HE3	1:C:93:MET:HA	1.63	0.78
1:A:89:SER:HA	1:A:92:MET:HE3	1.64	0.77
1:C:33:TRP:CZ2	1:C:47:VAL:HG13	2.19	0.77
1:C:66:LEU:HD11	1:C:172:TRP:O	1.85	0.77
1:C:97:GLY:O	1:C:288:ILE:HD11	1.86	0.76
1:A:8:ARG:HD3	1:A:16:TYR:CE2	2.21	0.76
1:A:44:GLN:HA	1:A:44:GLN:HE21	1.52	0.74
1:B:114:VAL:HG13	1:B:141:MET:HE1	1.69	0.74
1:B:116:GLN:HG3	1:B:199:LEU:HD13	1.70	0.74
1:A:29:MET:CE	1:A:31:GLU:HG2	2.10	0.73
1:B:39:GLU:HG3	1:B:44:GLN:NE2	2.04	0.72
1:A:124:ARG:HH12	1:A:182:LYS:HB2	1.53	0.72
1:B:8:ARG:HH11	1:B:8:ARG:N	1.87	0.72
1:C:192:ASN:N	1:C:192:ASN:HD22	1.88	0.72
1:A:276:GLU:O	1:A:280:THR:HG22	1.89	0.72
1:A:207:GLN:HB3	3:A:600:EDO:H22	1.72	0.71
1:A:7:VAL:CG1	1:A:78:PRO:HB2	2.21	0.71
1:B:305:THR:HB	1:B:306:PRO:CD	2.17	0.71
1:B:49:LYS:O	1:B:53:THR:HG23	1.91	0.70
1:A:174:THR:HG22	1:A:176:SER:O	1.91	0.70
1:A:93:MET:HE3	1:A:97:GLY:HA3	1.73	0.70
1:C:20:LYS:HG2	1:C:23:TYR:OH	1.92	0.70
1:B:93:MET:HE1	1:B:288:ILE:CG1	2.22	0.69
1:C:144:PHE:C	1:C:145:ASN:HD22	2.01	0.69
1:A:88:LYS:HG2	1:A:231:VAL:HG13	1.75	0.69
1:C:171:GLU:HB2	1:C:173:LYS:HE3	1.74	0.68
1:B:269:ASP:HB3	1:B:272:ILE:CG1	2.22	0.68
1:C:33:TRP:NE1	1:C:47:VAL:HA	2.08	0.68
1:C:88:LYS:HZ1	1:C:234:PHE:HB3	1.58	0.68
1:B:93:MET:CB	1:B:226:MET:HE1	2.23	0.68
1:B:288:ILE:C	1:B:288:ILE:HD13	2.19	0.68
1:B:114:VAL:CG1	1:B:141:MET:HE1	2.23	0.67
1:C:98:GLY:O	1:C:100:LYS:HG3	1.93	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:139:LEU:HB2	1:C:141:MET:HE2	1.74	0.67
1:B:106:HIS:HD2	1:B:200:TYR:OH	1.78	0.67
1:B:174:THR:HG22	1:B:176:SER:O	1.94	0.67
1:A:92:MET:HE1	1:A:231:VAL:HA	1.75	0.67
1:C:148:THR:CG2	1:C:152:GLN:HA	2.25	0.67
1:A:91:SER:O	1:A:95:MET:HG2	1.95	0.66
1:B:232:GLU:H	1:B:232:GLU:CD	2.01	0.66
1:C:91:SER:O	1:C:95:MET:HG2	1.93	0.66
1:C:198:ARG:HD2	1:C:221:HIS:HB3	1.76	0.66
1:A:8:ARG:O	1:A:13:ALA:HB2	1.96	0.66
1:B:194:ILE:N	1:B:194:ILE:HD12	2.11	0.65
1:C:136:THR:HG21	1:C:166:THR:OG1	1.97	0.65
1:A:29:MET:HE3	1:A:31:GLU:CG	2.12	0.65
1:C:204:LEU:HD12	1:C:224:TRP:CZ2	2.32	0.65
1:A:223:MET:HG2	4:A:940:HOH:O	1.97	0.65
1:C:88:LYS:NZ	1:C:234:PHE:HB3	2.12	0.65
1:A:114:VAL:HG13	1:A:141:MET:HE1	1.79	0.64
1:A:198:ARG:HD2	1:A:201:ARG:NH1	2.10	0.64
1:C:52:PHE:HZ	1:C:161:VAL:HG11	1.63	0.64
1:A:198:ARG:HD3	1:A:221:HIS:HB3	1.80	0.64
1:B:8:ARG:HD3	1:B:13:ALA:HA	1.80	0.64
1:B:93:MET:HE3	1:B:97:GLY:HA3	1.80	0.64
1:A:136:THR:CG2	1:A:166:THR:OG1	2.40	0.64
1:C:57:GLU:O	1:C:67:SER:HB3	1.97	0.64
1:A:281:LEU:HD11	1:A:288:ILE:HD11	1.79	0.63
1:C:219:GLY:HA3	1:C:223:MET:HE3	1.79	0.63
1:A:20:LYS:HA	1:A:23:TYR:CE2	2.33	0.63
1:B:194:ILE:HD12	1:B:194:ILE:H	1.64	0.62
1:C:145:ASN:HD22	1:C:145:ASN:N	1.98	0.62
1:A:124:ARG:NH1	1:A:182:LYS:HB2	2.14	0.62
1:C:268:VAL:HG12	1:C:269:ASP:N	2.15	0.61
1:C:95:MET:C	1:C:97:GLY:H	2.08	0.61
1:A:174:THR:CG2	1:A:176:SER:O	2.48	0.61
1:C:48:ASP:OD2	1:C:51:VAL:HG23	2.00	0.61
1:A:114:VAL:HG12	3:A:612:EDO:H12	1.81	0.61
1:B:70:GLN:HG2	1:B:75:ARG:NH1	2.16	0.60
1:C:127:THR:O	1:C:128:ASP:HB2	2.01	0.60
1:C:148:THR:HG23	1:C:152:GLN:HA	1.83	0.60
1:C:305:THR:HB	1:C:306:PRO:CD	2.31	0.60
1:C:77:ARG:HB2	1:C:164:ASN:ND2	2.17	0.60
1:C:288:ILE:HD13	1:C:288:ILE:C	2.27	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:8:ARG:HD3	1:B:13:ALA:CA	2.32	0.60
1:C:82:LEU:HB2	1:C:160:ALA:HB3	1.84	0.60
1:C:280:THR:O	1:C:284:THR:HG23	2.02	0.59
1:A:50:ASP:HB3	1:A:54:ARG:HH21	1.67	0.59
1:B:70:GLN:O	1:B:71:ASP:HB2	2.03	0.59
1:C:29:MET:HG2	1:C:31:GLU:HB3	1.84	0.59
1:B:298:ARG:O	1:B:302:ARG:HG2	2.04	0.58
1:C:54:ARG:NH1	1:C:61:PRO:HA	2.19	0.58
1:A:26:LEU:HG	1:A:147:ASP:HB3	1.84	0.58
1:C:87:PRO:HG2	1:C:90:VAL:CG2	2.34	0.58
1:B:269:ASP:CB	1:B:272:ILE:HG12	2.33	0.57
1:B:29:MET:HB3	1:B:31:GLU:HG3	1.85	0.57
1:B:106:HIS:CD2	1:B:200:TYR:OH	2.57	0.57
1:B:2:MET:HA	1:B:83:THR:O	2.05	0.57
1:A:127:THR:O	1:A:130:GLN:HG2	2.05	0.57
1:B:124:ARG:HH22	1:B:182:LYS:HB2	1.70	0.56
1:C:93:MET:HE1	1:C:288:ILE:HG12	1.86	0.56
1:C:96:LEU:C	1:C:96:LEU:HD12	2.30	0.56
1:C:130:GLN:O	1:C:130:GLN:HG3	2.06	0.56
1:C:92:MET:O	1:C:95:MET:O	2.24	0.56
1:B:29:MET:SD	1:B:31:GLU:HG2	2.45	0.56
1:B:104:ASP:O	1:B:108:GLN:HG3	2.06	0.56
1:B:124:ARG:NH2	1:B:182:LYS:HB2	2.20	0.56
1:B:91:SER:HB3	1:B:95:MET:HE3	1.87	0.55
1:C:25:VAL:HG21	1:C:277:TRP:HZ3	1.71	0.55
1:A:20:LYS:HA	1:A:23:TYR:CZ	2.41	0.55
1:C:127:THR:O	1:C:130:GLN:HG2	2.07	0.55
1:C:139:LEU:CB	1:C:141:MET:HE2	2.36	0.55
1:A:24:TYR:CD2	1:A:270:PRO:HB2	2.41	0.55
1:C:90:VAL:HA	1:C:226:MET:HE2	1.87	0.55
1:C:210:ALA:C	1:C:212:GLY:H	2.13	0.55
1:C:80:TYR:HE1	1:C:185:PHE:HD2	1.53	0.55
1:A:106:HIS:HD2	1:A:200:TYR:OH	1.89	0.55
1:C:29:MET:HE3	1:C:31:GLU:CG	2.16	0.55
1:C:292:ARG:O	1:C:295:ALA:HB3	2.07	0.55
1:C:204:LEU:HD13	1:C:204:LEU:O	2.06	0.55
1:A:50:ASP:HB3	1:A:54:ARG:NH2	2.22	0.55
1:C:20:LYS:HA	1:C:23:TYR:CZ	2.42	0.54
1:C:33:TRP:CD1	1:C:47:VAL:HG22	2.43	0.54
1:A:7:VAL:HG13	1:A:78:PRO:HB2	1.88	0.54
1:A:300:ASP:CG	1:B:36:ARG:HH22	2.15	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:33:TRP:CZ3	1:C:43:LEU:HD13	2.42	0.54
1:C:53:THR:O	1:C:57:GLU:HG3	2.08	0.54
1:C:218:VAL:HG12	1:C:218:VAL:O	2.08	0.54
1:C:8:ARG:O	1:C:13:ALA:HB2	2.08	0.54
1:C:288:ILE:HG23	1:C:289:ARG:N	2.22	0.54
1:B:108:GLN:HB2	1:B:207:GLN:HE22	1.72	0.53
1:C:144:PHE:O	1:C:158:THR:HA	2.08	0.53
1:B:8:ARG:H	1:B:8:ARG:NH1	1.94	0.53
1:C:30:GLY:O	1:C:32:ARG:HG3	2.09	0.53
1:C:191:ALA:C	1:C:192:ASN:HD22	2.17	0.53
1:C:204:LEU:HD12	1:C:224:TRP:CH2	2.44	0.53
1:C:184:GLY:O	1:C:188:ASN:ND2	2.41	0.53
1:A:80:TYR:CD2	1:A:186:ILE:HD12	2.43	0.53
1:A:301:LEU:O	1:A:305:THR:OG1	2.23	0.53
1:C:220:LYS:O	1:C:223:MET:HG3	2.09	0.53
1:A:183:THR:HA	1:A:187:GLU:OE1	2.10	0.52
1:B:302:ARG:O	1:B:305:THR:HG23	2.09	0.52
1:C:95:MET:O	1:C:97:GLY:N	2.42	0.52
1:C:278:MET:O	1:C:282:LYS:HG3	2.10	0.52
1:B:143:LEU:CD2	1:B:160:ALA:HA	2.40	0.52
1:C:80:TYR:CD1	1:C:186:ILE:HD13	2.45	0.52
1:A:70:GLN:HB3	1:A:75:ARG:HG3	1.91	0.51
1:B:91:SER:HB3	1:B:95:MET:CE	2.41	0.51
1:C:9:SER:O	1:C:11:GLY:N	2.43	0.51
1:C:52:PHE:O	1:C:55:LEU:HB3	2.11	0.51
1:C:95:MET:C	1:C:97:GLY:N	2.69	0.51
1:B:30:GLY:O	1:B:32:ARG:HD2	2.11	0.51
1:A:92:MET:O	1:A:96:LEU:HB2	2.11	0.51
1:C:195:ALA:O	1:C:198:ARG:HB2	2.10	0.51
1:B:194:ILE:H	1:B:194:ILE:CD1	2.24	0.51
1:C:25:VAL:HG12	1:C:95:MET:HB2	1.93	0.51
1:C:33:TRP:NE1	1:C:47:VAL:HG22	2.26	0.50
1:B:8:ARG:H	1:B:8:ARG:HD2	1.75	0.50
1:C:270:PRO:O	1:C:273:LYS:HB3	2.10	0.50
1:A:122:SER:HA	1:A:136:THR:HB	1.92	0.50
1:A:25:VAL:HG21	1:A:277:TRP:HZ3	1.76	0.50
1:C:90:VAL:HA	1:C:226:MET:CE	2.42	0.50
1:C:138:ASN:O	1:C:139:LEU:HG	2.11	0.50
1:C:202:GLU:O	1:C:205:LYS:HB3	2.11	0.50
1:A:25:VAL:HG21	1:A:277:TRP:CZ3	2.47	0.50
1:C:55:LEU:HD21	1:C:140:VAL:CG1	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:108:GLN:OE1	1:A:207:GLN:NE2	2.44	0.50
1:C:7:VAL:HG23	1:C:80:TYR:C	2.37	0.50
1:C:112:PHE:O	1:C:116:GLN:HG2	2.12	0.50
1:C:114:VAL:HG13	1:C:141:MET:SD	2.52	0.50
1:A:136:THR:HG23	1:A:138:ASN:ND2	2.27	0.49
1:B:22:ASN:HB3	1:B:146:HIS:CD2	2.47	0.49
1:C:108:GLN:HB2	1:C:207:GLN:HE22	1.76	0.49
1:A:234:PHE:O	1:A:235:SER:HB3	2.12	0.49
1:C:96:LEU:HD13	1:C:281:LEU:HD12	1.93	0.49
1:A:41:LEU:HD13	1:A:61:PRO:HD2	1.93	0.49
1:A:286:PHE:HD1	1:A:288:ILE:HD12	1.76	0.49
1:C:278:MET:HB3	1:C:282:LYS:HE3	1.95	0.49
1:C:88:LYS:HG3	1:C:92:MET:HE2	1.94	0.49
1:A:298:ARG:O	1:A:302:ARG:HG3	2.13	0.49
1:B:19:ASP:HB3	1:B:22:ASN:ND2	2.28	0.49
1:B:33:TRP:CZ2	1:B:47:VAL:HG13	2.48	0.49
1:A:1:MET:HB3	1:A:201:ARG:HH21	1.78	0.49
1:C:164:ASN:OD1	1:C:174:THR:HG23	2.13	0.49
1:A:1:MET:HB3	1:A:201:ARG:NH2	2.28	0.49
1:B:93:MET:HB2	1:B:226:MET:CE	2.40	0.49
1:B:69:MET:SD	1:B:74:ASN:HB2	2.53	0.48
1:C:210:ALA:C	1:C:212:GLY:N	2.70	0.48
1:C:204:LEU:HD12	1:C:224:TRP:HZ2	1.77	0.48
1:A:92:MET:CE	1:A:231:VAL:HG22	2.43	0.48
1:B:68:ARG:O	1:B:75:ARG:HG2	2.14	0.48
1:B:300:ASP:O	1:B:304:LEU:HG	2.13	0.48
1:C:287:ASP:HB3	1:C:290:ALA:HB3	1.96	0.48
1:C:51:VAL:C	1:C:54:ARG:HG2	2.38	0.48
1:A:93:MET:HE1	1:A:288:ILE:HG23	1.93	0.48
1:C:281:LEU:HD23	1:C:281:LEU:O	2.14	0.48
1:A:54:ARG:HD2	4:A:957:HOH:O	2.12	0.48
1:C:124:ARG:HH22	1:C:182:LYS:HB2	1.78	0.48
1:A:8:ARG:HD3	1:A:16:TYR:CZ	2.49	0.48
1:B:10:ALA:CB	1:B:57:GLU:HG3	2.44	0.48
1:C:12:SER:C	1:C:14:GLY:H	2.21	0.48
1:A:125:VAL:HG13	1:A:125:VAL:O	2.14	0.47
1:C:77:ARG:O	1:C:164:ASN:HB3	2.14	0.47
1:C:100:LYS:HB3	1:C:103:ILE:HD12	1.95	0.47
1:C:120:LEU:O	1:C:121:ALA:C	2.56	0.47
1:C:9:SER:C	1:C:11:GLY:N	2.72	0.47
1:C:23:TYR:HB3	1:C:145:ASN:O	2.12	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:80:TYR:HE1	1:C:185:PHE:CD2	2.31	0.47
1:C:107:ASN:HA	4:C:910:HOH:O	2.13	0.47
1:A:92:MET:HE1	1:A:231:VAL:HG22	1.97	0.47
1:C:126:MET:HA	1:C:131:SER:HA	1.96	0.47
1:A:229:VAL:HA	1:A:230:PRO:HD3	1.71	0.47
1:C:161:VAL:HG12	4:C:906:HOH:O	2.14	0.47
1:C:294:ALA:O	1:C:298:ARG:HB2	2.14	0.47
1:B:93:MET:CG	1:B:226:MET:HE1	2.44	0.47
1:C:23:TYR:CD2	1:C:29:MET:HB2	2.50	0.47
1:C:60:LEU:HB3	1:C:61:PRO:HD2	1.96	0.47
1:C:117:VAL:HG22	1:C:196:PHE:CE1	2.49	0.47
1:C:142:ALA:O	1:C:144:PHE:CE1	2.68	0.47
1:B:32:ARG:NH2	1:B:111:ASP:OD1	2.47	0.47
1:C:185:PHE:O	1:C:188:ASN:N	2.47	0.46
1:A:297:GLN:O	1:A:301:LEU:HB2	2.15	0.46
1:C:55:LEU:HD21	1:C:140:VAL:HG12	1.97	0.46
1:C:192:ASN:N	1:C:192:ASN:ND2	2.58	0.46
1:B:12:SER:O	1:B:13:ALA:C	2.58	0.46
1:C:164:ASN:ND2	1:C:174:THR:HG23	2.30	0.46
1:B:205:LYS:O	1:B:209:GLU:HG3	2.14	0.46
1:B:8:ARG:CG	1:B:9:SER:H	2.24	0.46
1:B:48:ASP:OD2	1:B:51:VAL:HG23	2.15	0.46
1:C:204:LEU:HD13	1:C:204:LEU:C	2.40	0.46
1:A:286:PHE:CD1	1:A:288:ILE:HD12	2.51	0.46
1:B:93:MET:HG2	1:B:226:MET:HE1	1.97	0.45
1:C:87:PRO:HB2	1:C:225:GLU:OE1	2.15	0.45
1:C:29:MET:HG2	1:C:31:GLU:CB	2.46	0.45
1:A:95:MET:SD	1:A:147:ASP:HB2	2.57	0.45
1:B:288:ILE:C	1:B:288:ILE:CD1	2.90	0.45
1:B:302:ARG:HG2	1:B:302:ARG:H	1.61	0.45
1:C:7:VAL:HG23	1:C:79:GLY:C	2.42	0.45
1:C:174:THR:HG22	1:C:176:SER:N	2.31	0.45
1:A:32:ARG:HD2	1:A:143:LEU:HB2	1.98	0.45
1:A:39:GLU:OE2	1:B:292:ARG:NH1	2.49	0.45
1:B:174:THR:CG2	1:B:176:SER:O	2.62	0.45
1:A:8:ARG:O	1:A:9:SER:C	2.59	0.45
1:B:143:LEU:HD22	1:B:160:ALA:HA	1.98	0.45
1:B:145:ASN:ND2	1:B:158:THR:OG1	2.50	0.45
1:B:33:TRP:CE2	1:B:47:VAL:HG22	2.52	0.45
1:B:93:MET:HE1	1:B:288:ILE:CD1	2.47	0.45
1:C:2:MET:HE3	1:C:4:ILE:HG12	1.99	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:31:GLU:H	1:A:31:GLU:HG3	1.59	0.45
1:A:106:HIS:HE1	1:A:158:THR:OG1	2.00	0.45
1:C:305:THR:HB	1:C:306:PRO:HD3	1.98	0.45
1:A:32:ARG:NE	1:A:143:LEU:HD12	2.32	0.45
1:B:118:GLU:HA	1:B:139:LEU:CD1	2.47	0.44
1:C:118:GLU:HA	1:C:139:LEU:CD1	2.47	0.44
1:C:141:MET:HG2	1:C:162:VAL:HA	1.99	0.44
1:C:11:GLY:C	1:C:13:ALA:H	2.25	0.44
1:A:33:TRP:CE2	1:A:47:VAL:HG22	2.52	0.44
1:A:128:ASP:C	1:A:130:GLN:H	2.26	0.44
1:A:232:GLU:HG2	1:A:233:ALA:N	2.32	0.44
4:A:969:HOH:O	1:B:115:ARG:HD2	2.18	0.44
1:C:25:VAL:HG21	1:C:277:TRP:CZ3	2.52	0.44
1:C:56:LEU:C	1:C:58:GLY:H	2.24	0.44
1:B:226:MET:HE3	1:B:229:VAL:HG21	1.99	0.44
1:C:155:GLN:O	1:C:157:HIS:ND1	2.51	0.44
1:A:174:THR:HG23	1:A:176:SER:N	2.32	0.44
1:A:300:ASP:OD2	1:B:36:ARG:NH2	2.50	0.44
1:C:8:ARG:O	1:C:9:SER:C	2.60	0.44
1:C:23:TYR:CG	1:C:29:MET:HB2	2.53	0.44
1:C:36:ARG:O	1:C:39:GLU:HB3	2.18	0.44
1:C:36:ARG:HB2	1:C:138:ASN:CB	2.47	0.44
1:C:111:ASP:O	1:C:115:ARG:HG3	2.17	0.44
1:A:8:ARG:NH1	1:A:16:TYR:CE1	2.86	0.44
1:C:106:HIS:HE1	1:C:158:THR:OG1	2.01	0.44
1:C:268:VAL:HG12	1:C:272:ILE:HG13	1.99	0.44
1:C:281:LEU:HD23	1:C:281:LEU:C	2.42	0.44
1:A:127:THR:O	1:A:128:ASP:HB2	2.18	0.44
1:B:276:GLU:O	1:B:280:THR:HG22	2.17	0.44
1:C:116:GLN:O	1:C:119:ALA:HB3	2.18	0.44
1:C:232:GLU:C	1:C:234:PHE:H	2.26	0.44
1:C:48:ASP:C	1:C:50:ASP:H	2.26	0.43
1:C:93:MET:HA	1:C:93:MET:CE	2.42	0.43
1:C:277:TRP:O	1:C:280:THR:HG22	2.18	0.43
1:A:99:ASP:OD1	1:A:99:ASP:C	2.62	0.43
1:A:88:LYS:CG	1:A:231:VAL:HG13	2.48	0.43
1:C:174:THR:CG2	1:C:176:SER:H	2.32	0.43
1:B:214:GLU:O	1:B:227:PRO:HD3	2.18	0.43
1:C:25:VAL:CG1	1:C:95:MET:HB2	2.49	0.43
1:C:229:VAL:O	1:C:231:VAL:HG23	2.18	0.43
1:B:174:THR:HG23	1:B:176:SER:H	1.84	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:93:MET:HB2	1:A:226:MET:HE1	2.01	0.43
1:C:72:GLY:O	1:C:73:SER:HB3	2.19	0.43
1:C:269:ASP:HB3	1:C:272:ILE:HG12	2.01	0.43
1:C:276:GLU:O	1:C:280:THR:HG22	2.19	0.43
1:C:127:THR:O	1:C:128:ASP:CB	2.67	0.42
1:C:162:VAL:O	1:C:163:ALA:C	2.61	0.42
1:C:174:THR:CG2	1:C:175:LEU:N	2.82	0.42
1:B:10:ALA:HB1	1:B:57:GLU:HG3	2.00	0.42
1:B:141:MET:HG2	1:B:162:VAL:HG22	2.01	0.42
1:C:125:VAL:O	1:C:125:VAL:HG13	2.17	0.42
1:C:216:GLU:O	1:C:218:VAL:HG23	2.19	0.42
1:B:205:LYS:HD2	1:B:215:THR:OG1	2.19	0.42
1:A:203:LYS:O	1:A:207:GLN:HG2	2.19	0.42
1:B:48:ASP:OD2	1:B:50:ASP:HB2	2.19	0.42
1:B:124:ARG:HH22	1:B:182:LYS:CG	2.32	0.42
1:C:54:ARG:HE	1:C:54:ARG:HB2	1.63	0.42
1:C:164:ASN:CG	1:C:174:THR:HG23	2.44	0.42
1:C:51:VAL:O	1:C:54:ARG:HG2	2.19	0.42
1:C:145:ASN:N	1:C:145:ASN:ND2	2.66	0.42
1:C:288:ILE:HG23	1:C:289:ARG:HG3	2.02	0.42
1:B:93:MET:HE3	1:B:93:MET:HA	2.01	0.42
1:C:164:ASN:HD21	1:C:174:THR:HG23	1.85	0.42
1:A:201:ARG:NE	4:A:939:HOH:O	2.48	0.42
1:B:8:ARG:HG3	1:B:12:SER:OG	2.20	0.42
1:B:89:SER:C	1:B:226:MET:HE2	2.44	0.42
1:B:26:LEU:HG	1:B:147:ASP:HB3	2.02	0.42
1:B:108:GLN:CD	1:B:207:GLN:NE2	2.78	0.42
1:C:166:THR:HG23	1:C:167:GLN:N	2.34	0.42
1:C:11:GLY:C	1:C:13:ALA:N	2.77	0.41
1:B:70:GLN:O	1:B:71:ASP:CB	2.68	0.41
1:C:143:LEU:N	1:C:143:LEU:HD23	2.35	0.41
1:A:146:HIS:HB2	1:A:157:HIS:NE2	2.35	0.41
1:A:198:ARG:CD	1:A:221:HIS:O	2.68	0.41
1:B:152:GLN:HE21	1:B:152:GLN:HB3	1.53	0.41
1:C:11:GLY:O	1:C:13:ALA:N	2.53	0.41
1:C:80:TYR:CG	1:C:186:ILE:HD13	2.56	0.41
1:C:96:LEU:HD13	1:C:281:LEU:CD1	2.50	0.41
1:B:108:GLN:CB	1:B:207:GLN:HE22	2.34	0.41
1:A:92:MET:HE1	1:A:231:VAL:CA	2.47	0.41
1:C:36:ARG:HB2	1:C:138:ASN:HB2	2.02	0.41
1:C:94:ALA:HA	1:C:99:ASP:O	2.20	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:126:MET:HE2	1:A:129:GLY:C	2.46	0.41
1:B:124:ARG:HH22	1:B:182:LYS:CB	2.33	0.41
1:C:122:SER:HB3	1:C:135:LEU:HA	2.02	0.41
1:C:164:ASN:HD21	1:C:174:THR:CG2	2.33	0.41
1:B:168:HIS:O	1:B:169:ASN:C	2.64	0.41
1:C:23:TYR:CE2	1:C:29:MET:SD	3.13	0.41
1:C:36:ARG:HD3	1:C:36:ARG:H	1.86	0.41
1:C:105:ALA:HA	1:C:207:GLN:NE2	2.35	0.41
1:B:124:ARG:HH22	1:B:182:LYS:HD2	1.86	0.41
1:C:106:HIS:HA	1:C:204:LEU:HD23	2.04	0.40
1:C:272:ILE:HG22	1:C:276:GLU:OE1	2.22	0.40
1:A:121:ALA:CB	1:A:136:THR:HG22	2.51	0.40
1:C:148:THR:HA	1:C:153:GLU:O	2.21	0.40
1:A:121:ALA:HB1	1:A:136:THR:HG22	2.03	0.40
1:A:198:ARG:HD3	1:A:221:HIS:O	2.21	0.40
1:C:99:ASP:OD1	1:C:101:ARG:HB2	2.22	0.40
1:C:118:GLU:HA	1:C:139:LEU:HD12	2.03	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	272/330 (82%)	256 (94%)	13 (5%)	3 (1%)	11	25
1	B	272/330 (82%)	251 (92%)	16 (6%)	5 (2%)	6	14
1	C	272/330 (82%)	217 (80%)	39 (14%)	16 (6%)	1	1
All	All	816/990 (82%)	724 (89%)	68 (8%)	24 (3%)	3	6

All (24) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	9	SER
1	C	96	LEU
1	B	180	VAL
1	B	306	PRO
1	C	10	ALA
1	C	94	ALA
1	C	99	ASP
1	C	121	ALA
1	C	221	HIS
1	C	305	THR
1	B	71	ASP
1	B	182	LYS
1	C	12	SER
1	C	11	GLY
1	C	185	PHE
1	A	20	LYS
1	C	23	TYR
1	C	24	TYR
1	C	78	PRO
1	C	128	ASP
1	A	230	PRO
1	B	230	PRO
1	C	230	PRO
1	C	268	VAL

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	225/267 (84%)	206 (92%)	19 (8%)	10	23
1	B	225/267 (84%)	209 (93%)	16 (7%)	13	31
1	C	225/267 (84%)	215 (96%)	10 (4%)	25	50
All	All	675/801 (84%)	630 (93%)	45 (7%)	15	33

All (45) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	19	ASP
1	A	31	GLU
1	A	32	ARG
1	A	36	ARG
1	A	44	GLN
1	A	70	GLN
1	A	101	ARG
1	A	114	VAL
1	A	124	ARG
1	A	136	THR
1	A	174	THR
1	A	178	ASP
1	A	180	VAL
1	A	187	GLU
1	A	205	LYS
1	A	223	MET
1	A	280	THR
1	A	281	LEU
1	A	302	ARG
1	B	8	ARG
1	B	31	GLU
1	B	32	ARG
1	B	36	ARG
1	B	88	LYS
1	B	114	VAL
1	B	152	GLN
1	B	174	THR
1	B	203	LYS
1	B	204	LEU
1	B	218	VAL
1	B	267	HIS
1	B	280	THR
1	B	281	LEU
1	B	288	ILE
1	B	302	ARG
1	C	1	MET
1	C	32	ARG
1	C	96	LEU
1	C	143	LEU
1	C	145	ASN
1	C	174	THR
1	C	192	ASN
1	C	232	GLU

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Mol	Chain	Res	Type
1	C	270	PRO
1	C	288	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (38) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	6	GLN
1	A	15	ASN
1	A	44	GLN
1	A	70	GLN
1	A	76	HIS
1	A	106	HIS
1	A	107	ASN
1	A	130	GLN
1	A	145	ASN
1	A	152	GLN
1	A	297	GLN
1	B	6	GLN
1	B	70	GLN
1	B	106	HIS
1	B	107	ASN
1	B	130	GLN
1	B	145	ASN
1	B	152	GLN
1	B	193	GLN
1	B	207	GLN
1	B	221	HIS
1	B	297	GLN
1	C	6	GLN
1	C	22	ASN
1	C	70	GLN
1	C	106	HIS
1	C	107	ASN
1	C	130	GLN
1	C	138	ASN
1	C	145	ASN
1	C	146	HIS
1	C	152	GLN
1	C	169	ASN
1	C	188	ASN
1	C	192	ASN
1	C	207	GLN

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Mol	Chain	Res	Type
1	C	221	HIS
1	C	297	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 ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

Of 14 ligands modelled in this entry, 3 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	$\# Z > 2$
3	EDO	B	604	-	3,3,3	0.57	0	2,2,2	0.25	0
3	EDO	C	603	-	3,3,3	0.59	0	2,2,2	0.18	0
3	EDO	B	607	-	3,3,3	0.57	0	2,2,2	0.26	0
3	EDO	A	612	-	3,3,3	0.61	0	2,2,2	0.15	0
3	EDO	A	600	-	3,3,3	0.56	0	2,2,2	0.14	0
3	EDO	A	606	-	3,3,3	0.65	0	2,2,2	0.22	0
3	EDO	A	608	-	3,3,3	0.59	0	2,2,2	0.25	0
3	EDO	B	602	-	3,3,3	0.62	0	2,2,2	0.22	0
3	EDO	B	610	-	3,3,3	0.58	0	2,2,2	0.28	0
3	EDO	B	605	-	3,3,3	0.58	0	2,2,2	0.22	0
3	EDO	A	601	-	3,3,3	0.58	0	2,2,2	0.27	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
3	EDO	B	604	-	-	1/1/1/1	-
3	EDO	C	603	-	-	1/1/1/1	-
3	EDO	B	607	-	-	1/1/1/1	-
3	EDO	A	612	-	-	1/1/1/1	-
3	EDO	A	600	-	-	1/1/1/1	-
3	EDO	A	606	-	-	1/1/1/1	-
3	EDO	A	608	-	-	1/1/1/1	-
3	EDO	B	602	-	-	1/1/1/1	-
3	EDO	B	610	-	-	1/1/1/1	-
3	EDO	B	605	-	-	1/1/1/1	-
3	EDO	A	601	-	-	1/1/1/1	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (11) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	600	EDO	O1-C1-C2-O2
3	A	601	EDO	O1-C1-C2-O2
3	A	606	EDO	O1-C1-C2-O2
3	A	608	EDO	O1-C1-C2-O2
3	A	612	EDO	O1-C1-C2-O2
3	B	602	EDO	O1-C1-C2-O2
3	B	604	EDO	O1-C1-C2-O2
3	B	605	EDO	O1-C1-C2-O2
3	B	607	EDO	O1-C1-C2-O2
3	B	610	EDO	O1-C1-C2-O2
3	C	603	EDO	O1-C1-C2-O2

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	612	EDO	2	0
3	A	600	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

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	275/330 (83%)	-0.06	5 (1%) 67 63	18, 45, 73, 104	1 (0%)
1	B	276/330 (83%)	0.09	12 (4%) 40 34	29, 49, 84, 116	0
1	C	276/330 (83%)	1.53	82 (29%) 1 1	56, 105, 152, 161	0
All	All	827/990 (83%)	0.52	99 (11%) 9 7	18, 59, 139, 161	1 (0%)

All (99) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	167	GLN	5.0
1	C	26	LEU	4.9
1	C	134	VAL	4.8
1	C	172	TRP	4.4
1	C	17	TYR	4.3
1	C	125	VAL	4.3
1	C	131	SER	4.2
1	C	73	SER	4.1
1	C	127	THR	4.1
1	C	15	ASN	4.1
1	C	268	VAL	4.1
1	C	66	LEU	4.1
1	C	18	THR	4.0
1	C	76	HIS	3.8
1	C	166	THR	3.8
1	C	270	PRO	3.7
1	C	56	LEU	3.7
1	B	304	LEU	3.6
1	C	135	LEU	3.6
1	C	24	TYR	3.6
1	C	25	VAL	3.5
1	C	16	TYR	3.4
1	C	13	ALA	3.3

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Mol	Chain	Res	Type	RSRZ
1	C	51	VAL	3.2
1	C	55	LEU	3.2
1	C	38	ALA	3.2
1	B	180	VAL	3.2
1	C	64	ALA	3.1
1	C	272	ILE	3.1
1	C	29	MET	3.1
1	C	37	GLY	3.0
1	C	47	VAL	3.0
1	C	271	GLU	2.9
1	C	291	TYR	2.9
1	C	52	PHE	2.9
1	C	72	GLY	2.9
1	B	307	GLY	2.9
1	C	306	PRO	2.9
1	B	130	GLN	2.9
1	C	173	LYS	2.8
1	C	277	TRP	2.8
1	C	30	GLY	2.7
1	C	45	GLY	2.7
1	A	268	VAL	2.7
1	C	147	ASP	2.7
1	C	10	ALA	2.7
1	C	199	LEU	2.7
1	C	123	THR	2.7
1	C	233	ALA	2.7
1	C	65	ASP	2.7
1	C	41	LEU	2.6
1	B	75	ARG	2.6
1	C	33	TRP	2.6
1	C	53	THR	2.6
1	C	96	LEU	2.6
1	A	306	PRO	2.6
1	C	60	LEU	2.5
1	C	36	ARG	2.5
1	C	14	GLY	2.5
1	C	74	ASN	2.5
1	C	43	LEU	2.5
1	C	146	HIS	2.5
1	C	23	TYR	2.5
1	C	75	ARG	2.4
1	B	177	SER	2.4

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Mol	Chain	Res	Type	RSRZ
1	C	128	ASP	2.4
1	C	78	PRO	2.4
1	C	175	LEU	2.4
1	A	72	GLY	2.4
1	C	22	ASN	2.4
1	C	145	ASN	2.3
1	B	235	SER	2.3
1	C	21	ASP	2.3
1	B	288	ILE	2.3
1	C	129	GLY	2.3
1	C	69	MET	2.3
1	C	292	ARG	2.3
1	C	63	GLY	2.3
1	C	122	SER	2.2
1	B	306	PRO	2.2
1	C	137	GLY	2.2
1	C	282	LYS	2.2
1	C	267	HIS	2.2
1	C	305	THR	2.2
1	C	20	LYS	2.2
1	C	28	SER	2.2
1	C	274	MET	2.1
1	B	183	THR	2.1
1	B	70	GLN	2.1
1	C	177	SER	2.1
1	C	286	PHE	2.1
1	C	281	LEU	2.1
1	B	1	MET	2.1
1	A	307	GLY	2.1
1	C	161	VAL	2.1
1	C	62	ASP	2.0
1	C	126	MET	2.0
1	A	8	ARG	2.0
1	C	32	ARG	2.0

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

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
3	EDO	B	607	4/4	0.32	0.21	114,114,114,115	0
3	EDO	B	610	4/4	0.58	0.20	84,84,85,85	0
3	EDO	A	606	4/4	0.75	0.23	76,78,78,79	0
2	MG	C	902	1/1	0.79	0.19	78,78,78,78	0
3	EDO	A	608	4/4	0.81	0.14	75,78,78,78	0
3	EDO	A	601	4/4	0.82	0.17	84,85,86,87	0
3	EDO	B	602	4/4	0.83	0.15	74,76,77,79	0
3	EDO	B	604	4/4	0.84	0.11	86,86,86,86	0
3	EDO	B	605	4/4	0.84	0.22	68,69,70,70	0
3	EDO	C	603	4/4	0.87	0.17	89,89,89,90	0
3	EDO	A	612	4/4	0.89	0.24	57,59,59,61	0
3	EDO	A	600	4/4	0.90	0.17	63,64,65,65	0
2	MG	B	901	1/1	0.92	0.14	52,52,52,52	0
2	MG	A	900	1/1	0.99	0.03	36,36,36,36	0

6.5 Other polymers [i](#)

There are no such residues in this entry.