



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 9, 2026 – 06:30 AM UTC

PDB ID : 8XUB / pdb_00008xub
Title : Crystal structure of Y-50 TCR
Authors : Asa, M.; Nagae, M.; Yamasaki, S.
Deposited on : 2024-01-12
Resolution : 2.50 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

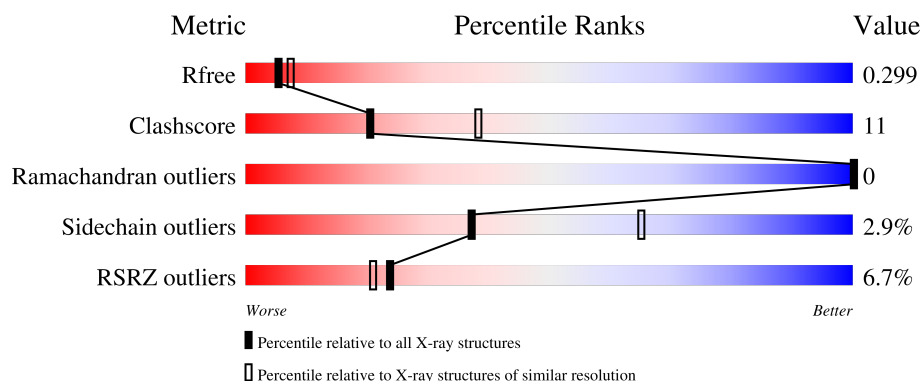
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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	5829 (2.50-2.50)
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

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	205	0% 85% 14% .
1	C	205	4% 70% 29% .
1	E	205	6% 69% 26% ..
1	G	205	13% 70% 28% .
2	B	249	2% 76% 22% .

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Mol	Chain	Length	Quality of chain
2	D	249	2%78%19%..
2	F	249	9%68%27%..
2	H	249	15%69%27%..

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 14371 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.

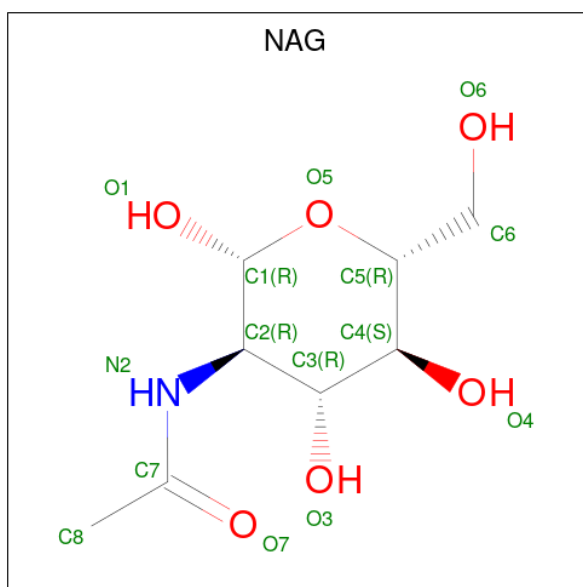
- Molecule 1 is a protein called Y-50 alpha.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	204	Total	C	N	O	S	0	0	0
			1616	1009	272	324	11			
1	C	204	Total	C	N	O	S	0	0	0
			1616	1009	272	324	11			
1	E	198	Total	C	N	O	S	0	0	0
			1569	979	265	314	11			
1	G	204	Total	C	N	O	S	0	0	0
			1616	1009	272	324	11			

- Molecule 2 is a protein called Y-50 beta.

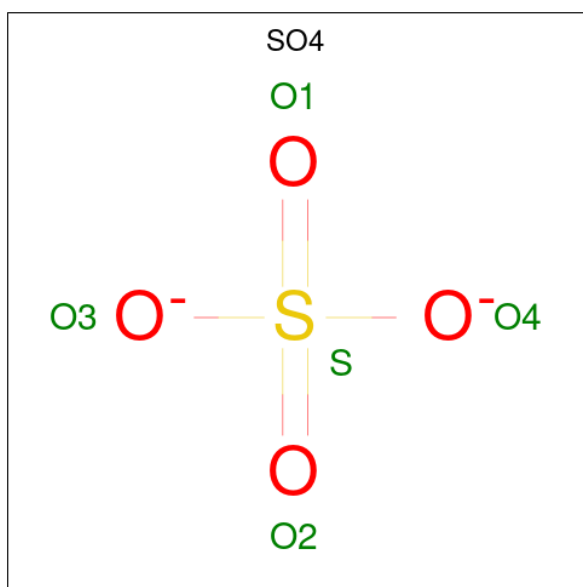
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	248	Total	C	N	O	S	0	0	0
			1974	1251	342	370	11			
2	D	243	Total	C	N	O	S	0	0	0
			1945	1232	337	365	11			
2	F	240	Total	C	N	O	S	0	0	0
			1929	1224	334	360	11			
2	H	240	Total	C	N	O	S	0	0	0
			1930	1224	334	361	11			

- Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (CCD ID: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			14	8	1	5		
3	B	1	Total	C	N	O	0	0
			14	8	1	5		
3	D	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	E	1	Total	C	N	O	0	0
			14	8	1	5		
3	F	1	Total	C	N	O	0	0
			14	8	1	5		
3	G	1	Total	C	N	O	0	0
			14	8	1	5		
3	H	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 4 is SULFATE ION (CCD ID: SO4) (formula: O₄S).



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

- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	4	Total	O	0	0
			4	4		
5	B	11	Total	O	0	0
			11	11		
5	C	5	Total	O	0	0
			5	5		
5	D	5	Total	O	0	0
			5	5		
5	E	2	Total	O	0	0
			2	2		
5	F	2	Total	O	0	0
			2	2		

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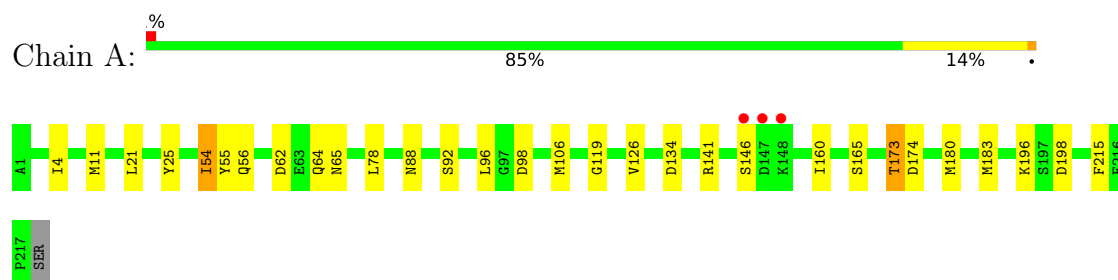
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Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	H	5	Total	O	0	0
			5	5		

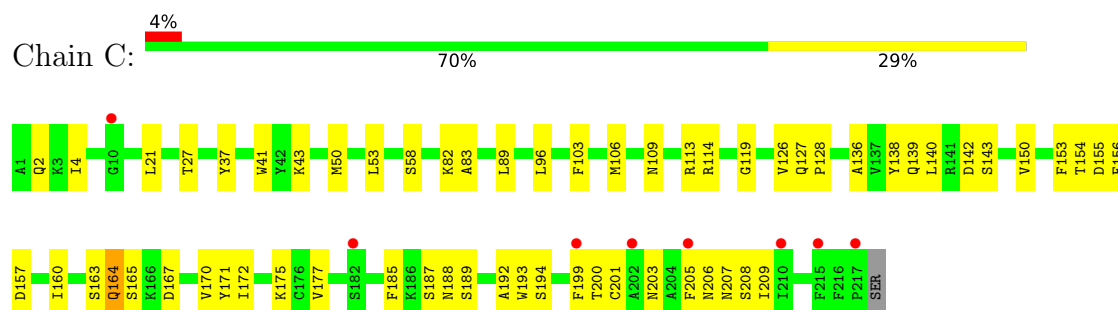
3 Residue-property plots [i](#)

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.

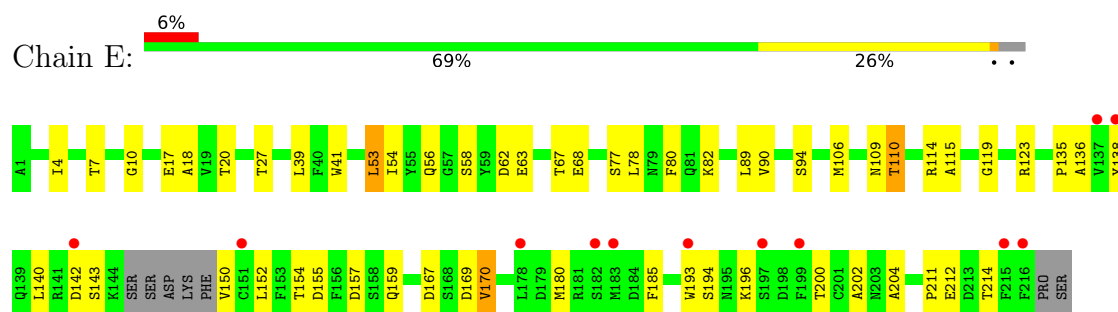
- Molecule 1: Y-50 alpha



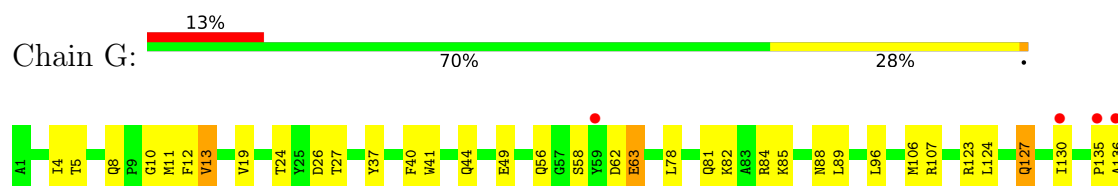
- Molecule 1: Y-50 alpha

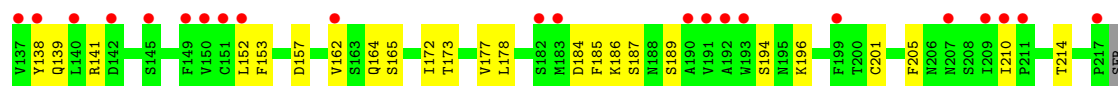


- Molecule 1: Y-50 alpha

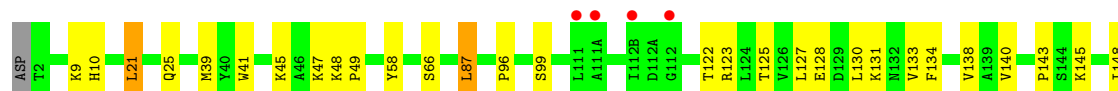
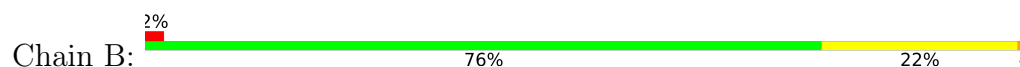


- Molecule 1: Y-50 alpha

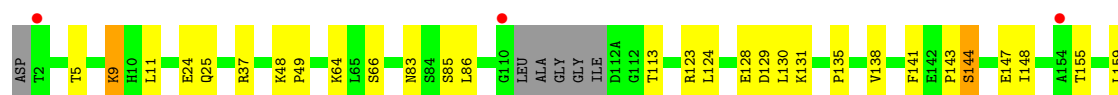
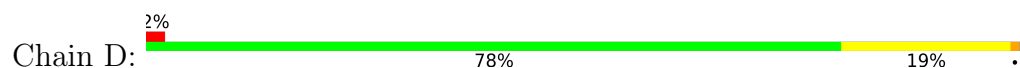




• Molecule 2: Y-50 beta



• Molecule 2: Y-50 beta

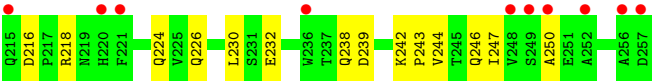


• Molecule 2: Y-50 beta



• Molecule 2: Y-50 beta





4 Data and refinement statistics

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants a, b, c, α , β , γ	68.11Å 168.69Å 173.12Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	44.03 – 2.50 44.03 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (44.03-2.50) 100.0 (44.03-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.74 (at 2.51Å)	Xtriage
Refinement program	PHENIX 1.20.1_4487	Depositor
R, R_{free}	0.226 , 0.299 0.226 , 0.299	Depositor DCC
R_{free} test set	3386 reflections (4.84%)	wwPDB-VP
Wilson B-factor (Å ²)	59.7	Xtriage
Anisotropy	0.326	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.31 , 40.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Estimated twinning fraction	0.015 for -h,l,k	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	14371	wwPDB-VP
Average B, all atoms (Å ²)	69.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.39% 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 [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: NAG, 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	0.45	0/1649	0.63	0/2228
1	C	0.40	0/1649	0.59	0/2228
1	E	0.35	0/1599	0.56	0/2159
1	G	0.38	0/1649	0.59	0/2228
2	B	0.45	0/2030	0.64	0/2761
2	D	0.42	0/2000	0.61	0/2719
2	F	0.37	0/1984	0.60	0/2698
2	H	0.39	0/1985	0.59	0/2699
All	All	0.40	0/14545	0.60	0/19720

There are no bond length outliers.

There are no bond angle outliers.

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	1616	0	1532	21	0
1	C	1616	0	1534	47	0
1	E	1569	0	1488	43	0
1	G	1616	0	1533	42	0
2	B	1974	0	1911	41	1

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	D	1945	0	1877	33	1
2	F	1929	0	1867	63	1
2	H	1930	0	1867	47	1
3	A	14	0	13	2	0
3	B	14	0	13	0	0
3	D	14	0	13	0	0
3	E	28	0	26	0	0
3	F	14	0	13	0	0
3	G	14	0	13	1	0
3	H	14	0	13	0	0
4	A	5	0	0	0	0
4	B	5	0	0	1	0
4	C	5	0	0	0	0
4	D	5	0	0	0	0
4	E	5	0	0	0	0
4	G	5	0	0	0	0
5	A	4	0	0	0	0
5	B	11	0	0	0	0
5	C	5	0	0	0	0
5	D	5	0	0	0	0
5	E	2	0	0	0	0
5	F	2	0	0	0	0
5	H	5	0	0	0	0
All	All	14371	0	13713	315	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 11.

All (315) 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:G:13:VAL:HG11	1:G:19:VAL:HG22	1.42	0.99
2:D:236:TRP:HB2	2:D:242:LYS:HG3	1.45	0.96
1:G:12:PHE:HD1	1:G:127:GLN:HE22	1.14	0.92
1:C:156:PHE:HB2	1:C:160:ILE:HD13	1.53	0.88
1:G:135:PRO:HB2	1:G:214:THR:HG22	1.57	0.86
1:G:130:ILE:HG21	1:G:157:ASP:HA	1.65	0.79
2:F:160:ALA:HB2	2:F:225:VAL:HG21	1.64	0.78
2:H:131:LYS:NZ	2:H:238:GLN:OE1	2.16	0.78
1:E:20:THR:HG22	1:E:90:VAL:HG22	1.66	0.77
1:C:43:LYS:HB3	1:C:53:LEU:HD21	1.67	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:163:SER:N	1:C:208:SER:OG	2.18	0.76
1:E:135:PRO:HB2	1:E:214:THR:HG22	1.70	0.73
2:H:6:GLN:HB2	2:H:120:PRO:HD2	1.71	0.73
1:G:139:GLN:NE2	1:G:201:CYS:SG	2.62	0.71
1:E:4:ILE:HD11	1:E:106:MET:HB2	1.70	0.71
1:G:210:ILE:HG22	1:G:214:THR:HG21	1.72	0.71
1:E:27:THR:HB	1:E:106:MET:HE1	1.72	0.70
1:G:26:ASP:OD1	1:G:85:LYS:NZ	2.25	0.70
1:E:54:ILE:HD12	1:E:68:GLU:HB2	1.76	0.68
2:B:127:LEU:HD21	2:B:164:TYR:HE1	1.59	0.68
2:F:147:GLU:OE1	2:F:155:THR:OG1	2.09	0.68
2:H:138:VAL:HG23	2:H:250:ALA:HB2	1.76	0.67
2:H:242:LYS:HG2	2:H:244:VAL:HG13	1.75	0.67
2:F:226:GLN:NE2	2:F:246:GLN:O	2.28	0.66
1:A:56:GLN:HB2	1:A:78:LEU:HD23	1.78	0.66
2:F:138:VAL:HG23	2:F:250:ALA:HB2	1.77	0.66
1:C:206:ASN:OD1	1:C:207:ASN:ND2	2.28	0.65
2:B:133:VAL:C	2:B:134:PHE:HD1	2.05	0.65
2:B:134:PHE:HD2	2:B:200:ARG:CZ	2.11	0.64
2:H:91:LEU:HD13	2:H:94:LEU:HD11	1.80	0.64
1:E:200:THR:HG23	1:E:202:ALA:H	1.63	0.63
1:C:138:TYR:CD1	2:D:147:GLU:HB2	2.33	0.63
1:C:153:PHE:HB2	1:C:205:PHE:CE2	2.33	0.63
2:H:101:LEU:HD13	2:H:123:ARG:HG3	1.80	0.63
1:G:139:GLN:OE1	1:G:141:ARG:NH2	2.31	0.63
2:B:156:LEU:HD13	2:B:207:LEU:HD23	1.81	0.63
2:F:193:GLN:CD	2:F:196:LEU:H	2.07	0.62
2:H:143:PRO:HD2	2:H:214:TRP:CZ2	2.35	0.62
2:H:138:VAL:HA	2:H:159:LEU:O	2.01	0.61
2:B:130:LEU:O	2:B:133:VAL:HG12	2.01	0.61
1:E:170:VAL:HA	1:E:194:SER:OG	2.01	0.61
1:C:150:VAL:HG12	1:C:193:TRP:HB3	1.83	0.60
1:G:81:GLN:HB3	1:G:84:ARG:HD2	1.85	0.59
1:C:139:GLN:O	2:D:144:SER:HB2	2.03	0.59
2:H:224:GLN:HE21	2:H:247:ILE:HD12	1.67	0.59
1:C:50:MET:HE1	1:C:103:PHE:CD2	2.37	0.59
2:H:2:THR:HG22	2:H:2:THR:O	2.03	0.58
2:F:25:GLN:O	2:F:85:SER:HB2	2.04	0.58
2:F:193:GLN:NE2	2:F:195:ALA:H	2.02	0.57
2:B:134:PHE:CD2	2:B:200:ARG:CZ	2.87	0.57
2:F:154:ALA:N	2:F:209:VAL:O	2.38	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:211:PRO:O	1:E:214:THR:HG23	2.05	0.56
2:D:129:ASP:OD2	2:D:131:LYS:HE3	2.05	0.56
2:D:143:PRO:HD2	2:D:214:TRP:CZ2	2.41	0.56
1:E:136:ALA:HA	1:E:214:THR:HB	1.87	0.56
1:A:25:TYR:CD1	1:A:106:MET:HE1	2.40	0.56
1:A:78:LEU:HD12	1:A:88:ASN:O	2.06	0.55
1:E:212:GLU:N	1:E:212:GLU:OE1	2.38	0.55
1:C:150:VAL:HG13	2:D:141:PHE:CE1	2.42	0.55
1:G:56:GLN:HB2	1:G:78:LEU:HD23	1.89	0.55
1:C:50:MET:HE1	1:C:103:PHE:CE2	2.42	0.55
2:H:193:GLN:HB3	2:H:196:LEU:HD23	1.87	0.55
2:H:226:GLN:HG2	2:H:247:ILE:CD1	2.37	0.55
1:C:200:THR:HG23	1:C:203:ASN:HB3	1.89	0.54
2:F:193:GLN:HE22	2:F:195:ALA:H	1.54	0.54
2:B:186:ASP:OD1	2:B:206:ARG:NH2	2.39	0.54
1:E:80:PHE:CZ	1:E:82:LYS:HG3	2.42	0.54
2:H:129:ASP:OD1	2:H:131:LYS:HB2	2.07	0.54
1:C:200:THR:CG2	1:C:203:ASN:HB3	2.36	0.54
1:A:96:LEU:HD23	1:A:126:VAL:HG12	1.90	0.54
2:H:209:VAL:HG12	2:H:210:SER:H	1.73	0.54
2:F:137:GLU:N	2:F:161:THR:O	2.39	0.54
1:A:65:ASN:HB2	3:A:301:NAG:N2	2.23	0.54
2:B:41:TRP:CD1	2:B:87:LEU:HD21	2.43	0.53
1:A:180:MET:HE2	1:A:183:MET:HG2	1.91	0.53
2:H:216:ASP:OD1	2:H:218:ARG:HG3	2.08	0.53
2:F:140:VAL:HG23	2:F:250:ALA:HB3	1.90	0.53
2:H:238:GLN:HG3	2:H:239:ASP:N	2.24	0.53
2:F:158:CYS:HB2	2:F:172:TRP:CH2	2.43	0.53
2:D:230:LEU:HD13	2:D:234:ASP:HB3	1.91	0.53
2:F:148:ILE:HG23	2:F:211:ALA:HB1	1.91	0.53
2:F:150:ARG:HB3	2:F:150:ARG:HH11	1.74	0.53
2:H:131:LYS:NZ	2:H:238:GLN:HB2	2.24	0.53
2:H:226:GLN:HG2	2:H:247:ILE:HD13	1.91	0.53
2:F:226:GLN:HE21	2:F:227:PHE:N	2.06	0.52
2:H:171:SER:OG	2:H:224:GLN:OE1	2.27	0.52
2:H:230:LEU:HD12	2:H:230:LEU:H	1.74	0.52
1:G:205:PHE:O	1:G:210:ILE:HD11	2.10	0.52
1:A:65:ASN:HB2	3:A:301:NAG:C7	2.40	0.52
2:D:236:TRP:CB	2:D:242:LYS:HG3	2.29	0.52
2:H:135:PRO:HD3	2:H:243:PRO:HB3	1.91	0.51
1:G:40:PHE:HE1	1:G:107:ARG:HG2	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:135:PRO:HD3	2:D:243:PRO:HB3	1.92	0.51
1:E:200:THR:O	1:E:204:ALA:N	2.42	0.51
2:F:162:GLY:O	2:F:200:ARG:NE	2.33	0.51
1:A:160:ILE:HD12	1:A:160:ILE:H	1.76	0.51
2:D:123:ARG:HD3	2:D:167:HIS:CG	2.46	0.51
1:E:109:ASN:O	1:E:114:ARG:NH1	2.43	0.51
1:C:127:GLN:OE1	1:C:175:LYS:NZ	2.37	0.51
1:A:54:ILE:HG13	1:A:55:TYR:N	2.25	0.51
1:E:150:VAL:HG12	1:E:193:TRP:HB3	1.92	0.50
2:F:155:THR:HG23	2:F:208:ARG:HG3	1.92	0.50
2:F:164:TYR:HB3	2:F:165:PRO:HD3	1.93	0.50
2:D:11:LEU:HD23	2:D:124:LEU:HD13	1.93	0.50
1:G:96:LEU:HD22	1:G:177:VAL:HG11	1.92	0.50
1:A:173:THR:HG22	1:A:174:ASP:O	2.12	0.50
1:E:17:GLU:HG3	1:E:18:ALA:H	1.77	0.50
2:D:25:GLN:O	2:D:85:SER:HB2	2.12	0.50
1:E:167:ASP:HB2	1:E:170:VAL:HG22	1.92	0.50
1:G:152:LEU:HD11	1:G:189:SER:HB2	1.94	0.50
2:B:25:GLN:HG2	2:B:39:MET:HE3	1.94	0.50
2:D:193:GLN:HB2	2:D:196:LEU:HD13	1.93	0.50
1:E:41:TRP:CE2	1:E:89:LEU:HB2	2.47	0.50
1:C:154:THR:OG1	1:C:155:ASP:N	2.45	0.49
2:H:166:PRO:HB3	2:H:201:TYR:CG	2.47	0.49
2:F:135:PRO:HB3	2:F:246:GLN:OE1	2.11	0.49
2:B:236:TRP:HB2	2:B:242:LYS:HD2	1.94	0.49
2:H:81:PRO:HG2	2:H:86:LEU:HD23	1.94	0.49
1:G:10:GLY:HA2	1:G:123:ARG:O	2.12	0.49
2:B:143:PRO:HD2	2:B:214:TRP:CZ2	2.47	0.49
1:E:10:GLY:HA2	1:E:123:ARG:O	2.12	0.49
2:F:193:GLN:OE1	2:F:196:LEU:N	2.41	0.49
1:C:140:LEU:HD21	2:D:155:THR:O	2.11	0.49
1:E:150:VAL:HG13	2:F:141:PHE:CE1	2.48	0.49
2:B:193:GLN:OE1	2:B:199:SER:HA	2.12	0.49
1:G:184:ASP:OD1	1:G:184:ASP:O	2.30	0.49
2:F:83:ASN:HB2	2:F:86:LEU:HD23	1.94	0.48
2:F:127:LEU:HD21	2:F:164:TYR:CE2	2.48	0.48
1:G:185:PHE:CE2	1:G:187:SER:HB3	2.48	0.48
2:F:150:ARG:HB3	2:F:150:ARG:NH1	2.28	0.48
1:G:136:ALA:HB3	1:G:138:TYR:CE1	2.48	0.48
1:A:141:ARG:HD2	1:A:146:SER:O	2.14	0.48
2:F:213:PHE:CE2	2:F:221:PHE:HE1	2.31	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:13:VAL:CG1	1:G:19:VAL:HG22	2.29	0.48
1:G:210:ILE:CG2	1:G:214:THR:HG21	2.42	0.48
1:C:4:ILE:HD11	1:C:106:MET:HB2	1.94	0.48
1:E:58:SER:O	1:E:82:LYS:HD2	2.14	0.48
1:G:210:ILE:HD12	1:G:210:ILE:N	2.29	0.48
2:H:148:ILE:HG23	2:H:211:ALA:HB1	1.95	0.48
1:C:170:VAL:HG21	1:C:199:PHE:CE1	2.48	0.48
1:A:4:ILE:HD11	1:A:106:MET:HB3	1.96	0.47
2:B:47:LYS:HB2	2:B:47:LYS:HE2	1.64	0.47
2:B:125:THR:HG21	2:B:165:PRO:HB3	1.96	0.47
1:E:56:GLN:HB2	1:E:78:LEU:HD21	1.95	0.47
2:H:242:LYS:HE3	2:H:244:VAL:CG1	2.44	0.47
2:F:154:ALA:O	2:F:208:ARG:HA	2.14	0.47
2:B:162:GLY:O	2:B:200:ARG:NH2	2.45	0.47
2:B:133:VAL:HG13	2:B:243:PRO:HG2	1.96	0.47
1:G:165:SER:HB2	1:G:172:ILE:HD11	1.96	0.47
1:C:113:ARG:CZ	2:D:66:SER:HB3	2.45	0.47
1:E:138:TYR:CD1	2:F:147:GLU:HB2	2.50	0.47
2:F:143:PRO:HD2	2:F:214:TRP:CH2	2.50	0.47
2:B:180:HIS:HB2	4:B:302:SO4:O1	2.15	0.47
1:G:58:SER:O	1:G:82:LYS:NZ	2.32	0.46
1:A:134:ASP:CG	2:B:150:ARG:HH22	2.24	0.46
1:C:2:GLN:HB3	1:C:106:MET:HE3	1.97	0.46
2:D:242:LYS:O	2:D:244:VAL:N	2.48	0.46
2:D:24:GLU:HG3	2:D:86:LEU:HD22	1.98	0.46
1:G:41:TRP:CE2	1:G:89:LEU:HB2	2.50	0.46
2:B:123:ARG:HD2	2:B:167:HIS:NE2	2.30	0.46
2:D:48:LYS:HA	2:D:48:LYS:HD3	1.71	0.46
1:A:98:ASP:HB2	1:A:126:VAL:HG21	1.97	0.46
2:B:215:GLN:HA	2:B:255:ARG:O	2.16	0.46
1:E:138:TYR:HB3	2:F:144:SER:OG	2.16	0.46
1:G:13:VAL:HG22	1:G:124:LEU:HD11	1.98	0.46
1:G:177:VAL:N	2:H:184:CYS:SG	2.89	0.46
2:B:47:LYS:O	2:B:48:LYS:HD3	2.16	0.46
2:D:217:PRO:HA	2:D:254:GLY:O	2.15	0.46
1:G:194:SER:OG	1:G:196:LYS:HG3	2.15	0.46
1:C:136:ALA:HB1	1:C:138:TYR:HE2	1.80	0.46
1:A:196:LYS:HE3	1:A:198:ASP:OD1	2.16	0.46
2:B:143:PRO:HD2	2:B:214:TRP:CE2	2.51	0.46
1:C:21:LEU:HD12	1:C:89:LEU:HD23	1.99	0.45
1:G:56:GLN:NE2	1:G:62:ASP:HB2	2.32	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:45:LYS:NZ	2:B:99:SER:O	2.49	0.45
1:C:170:VAL:HA	1:C:194:SER:HB3	1.97	0.45
1:C:185:PHE:CE2	1:C:187:SER:HB3	2.51	0.45
2:D:218:ARG:HE	2:D:218:ARG:HB3	1.39	0.45
2:B:9:LYS:HD3	2:B:10:HIS:NE2	2.32	0.45
2:B:21:LEU:HD22	2:B:122:THR:HG21	1.99	0.45
2:B:140:VAL:HG23	2:B:250:ALA:HB3	1.98	0.45
1:C:109:ASN:O	1:C:114:ARG:NH1	2.48	0.45
1:C:164:GLN:H	1:C:164:GLN:CD	2.11	0.45
2:F:136:PRO:HG3	2:F:227:PHE:HB2	1.99	0.45
1:C:150:VAL:HG12	1:C:193:TRP:CB	2.47	0.45
1:C:193:TRP:CD2	2:D:159:LEU:HD21	2.52	0.45
1:E:119:GLY:O	2:F:49:PRO:HB3	2.16	0.45
2:B:220:HIS:NE2	2:B:251:GLU:OE1	2.49	0.45
2:D:130:LEU:H	2:D:130:LEU:HD12	1.82	0.45
1:E:180:MET:HE3	2:F:208:ARG:O	2.17	0.45
1:C:4:ILE:HD11	1:C:106:MET:HE2	1.98	0.45
2:F:136:PRO:HD3	2:F:227:PHE:CD1	2.52	0.45
2:F:226:GLN:HE21	2:F:227:PHE:H	1.65	0.45
1:G:11:MET:HB3	1:G:11:MET:HE2	1.82	0.45
1:G:63:GLU:C	3:G:301:NAG:H81	2.42	0.45
2:H:21:LEU:HB2	2:H:89:LEU:HB3	1.99	0.45
2:B:242:LYS:HG2	2:B:244:VAL:HG13	1.99	0.44
2:D:236:TRP:CE2	2:D:243:PRO:HD3	2.52	0.44
1:G:4:ILE:HD11	1:G:106:MET:HB3	1.98	0.44
1:E:212:GLU:H	1:E:212:GLU:CD	2.24	0.44
2:F:191:LYS:HE2	2:F:199:SER:HB3	1.99	0.44
2:H:136:PRO:HB3	2:H:163:PHE:CD2	2.52	0.44
1:C:167:ASP:HB3	1:C:170:VAL:HG12	2.00	0.44
1:E:170:VAL:HG12	1:E:196:LYS:HD3	2.00	0.44
1:C:171:TYR:O	1:C:192:ALA:HA	2.18	0.44
2:F:220:HIS:CE1	2:F:222:ARG:HB2	2.53	0.44
2:H:25:GLN:O	2:H:85:SER:HB2	2.18	0.44
2:B:96:PRO:HG2	2:B:128:GLU:HG2	1.99	0.44
2:D:209:VAL:HG12	2:D:210:SER:N	2.33	0.44
2:F:172:TRP:CD2	2:F:207:LEU:HD22	2.52	0.44
2:H:173:TRP:NE1	2:H:224:GLN:OE1	2.46	0.44
1:E:62:ASP:O	1:E:63:GLU:HG2	2.18	0.44
1:E:115:ALA:HB1	2:F:52:LEU:HD22	1.98	0.44
1:E:138:TYR:HD1	2:F:144:SER:OG	2.01	0.44
2:F:183:VAL:HA	2:F:206:ARG:O	2.17	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:133:VAL:O	2:B:134:PHE:HD1	1.99	0.44
2:D:209:VAL:HG12	2:D:210:SER:H	1.82	0.44
1:E:185:PHE:HE1	2:F:208:ARG:HE	1.66	0.44
2:F:246:GLN:C	2:F:247:ILE:HD12	2.42	0.44
2:H:154:ALA:O	2:H:208:ARG:HA	2.18	0.44
1:C:136:ALA:HB1	1:C:138:TYR:CE2	2.53	0.43
1:G:96:LEU:HD11	1:G:186:LYS:HD3	2.00	0.43
1:A:215:PHE:CZ	2:B:150:ARG:HD3	2.52	0.43
1:A:119:GLY:O	2:B:49:PRO:HB3	2.18	0.43
2:B:25:GLN:CG	2:B:39:MET:HE3	2.48	0.43
2:B:145:LYS:H	2:B:145:LYS:HD2	1.83	0.43
2:D:232:GLU:HG3	2:D:233:ASN:N	2.33	0.43
2:B:148:ILE:HG23	2:B:211:ALA:HB1	2.00	0.43
2:F:129:ASP:HB3	2:F:131:LYS:HG3	2.00	0.43
1:G:13:VAL:HG21	1:G:19:VAL:CG2	2.48	0.43
1:G:164:GLN:OE1	1:G:164:GLN:HA	2.18	0.43
2:H:5:THR:HG22	2:H:24:GLU:HB2	1.99	0.43
2:B:166:PRO:HB3	2:B:201:TYR:CG	2.54	0.43
2:F:193:GLN:CD	2:F:196:LEU:HB2	2.44	0.43
2:H:181:ASP:OD1	2:H:182:GLY:N	2.51	0.43
2:D:9:LYS:HD3	2:D:9:LYS:N	2.33	0.43
2:D:209:VAL:HG11	2:D:213:PHE:CD2	2.53	0.43
2:F:220:HIS:CD2	2:F:251:GLU:HG3	2.54	0.43
1:G:27:THR:HG21	1:G:37:TYR:CD2	2.53	0.43
1:G:56:GLN:CB	1:G:78:LEU:HD23	2.49	0.43
2:H:135:PRO:HB3	2:H:246:GLN:NE2	2.33	0.43
1:C:41:TRP:CE2	1:C:89:LEU:HB2	2.54	0.43
1:C:164:GLN:OE1	1:C:164:GLN:N	2.39	0.43
1:E:41:TRP:O	1:E:53:LEU:HB2	2.19	0.43
2:F:172:TRP:CG	2:F:207:LEU:HD22	2.54	0.43
1:C:200:THR:OG1	1:C:201:CYS:N	2.52	0.43
1:E:136:ALA:HB3	1:E:138:TYR:HE2	1.83	0.43
1:E:140:LEU:HD13	1:E:152:LEU:HB3	2.00	0.43
1:E:169:ASP:HB3	1:E:196:LYS:HD2	2.00	0.43
2:F:193:GLN:NE2	2:F:196:LEU:H	2.17	0.43
1:A:11:MET:HG3	1:A:21:LEU:CD2	2.49	0.42
2:B:131:LYS:O	2:B:134:PHE:HE1	2.03	0.42
1:C:27:THR:HG21	1:C:37:TYR:CD2	2.54	0.42
2:D:195:ALA:C	2:D:196:LEU:HD12	2.44	0.42
2:H:145:LYS:HE3	2:H:145:LYS:HB2	1.70	0.42
2:F:89:LEU:HA	2:F:89:LEU:HD23	1.60	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:138:VAL:HG21	2:D:248:VAL:O	2.19	0.42
2:H:91:LEU:HB3	2:H:94:LEU:HD21	2.01	0.42
2:F:158:CYS:HB2	2:F:172:TRP:CZ2	2.53	0.42
1:G:78:LEU:HD12	1:G:88:ASN:O	2.20	0.42
1:E:154:THR:HG22	1:E:155:ASP:OD1	2.19	0.42
2:H:89:LEU:HD12	2:H:89:LEU:HA	1.80	0.42
1:C:58:SER:O	1:C:82:LYS:HE3	2.20	0.42
2:H:148:ILE:HG23	2:H:211:ALA:CB	2.50	0.42
1:C:4:ILE:CD1	1:C:106:MET:HE2	2.50	0.42
1:C:96:LEU:HA	1:C:126:VAL:HB	2.02	0.42
2:H:11:LEU:HD11	2:H:19:LYS:HE3	2.02	0.42
1:E:135:PRO:O	1:E:214:THR:HA	2.19	0.42
1:A:62:ASP:O	1:A:64:GLN:HG2	2.20	0.41
2:H:191:LYS:HG2	2:H:194:PRO:HA	2.02	0.41
1:C:128:PRO:HG3	1:C:177:VAL:HG13	2.01	0.41
1:G:44:GLN:HA	1:G:49:GLU:O	2.19	0.41
1:G:81:GLN:HG2	1:G:84:ARG:CZ	2.50	0.41
1:C:157:ASP:O	1:C:160:ILE:HD12	2.20	0.41
1:C:119:GLY:O	2:D:49:PRO:HB3	2.20	0.41
2:F:21:LEU:HD22	2:F:122:THR:HG21	2.02	0.41
2:F:48:LYS:HD3	2:F:48:LYS:HA	1.70	0.41
2:F:136:PRO:HB3	2:F:163:PHE:HB3	2.02	0.41
2:F:146:ALA:O	2:F:150:ARG:N	2.53	0.41
2:B:156:LEU:HD12	2:B:156:LEU:H	1.86	0.41
2:B:242:LYS:HA	2:B:243:PRO:HD3	1.86	0.41
2:F:137:GLU:O	2:F:161:THR:N	2.32	0.41
2:F:193:GLN:HA	2:F:194:PRO:HD3	1.85	0.41
2:H:137:GLU:O	2:H:161:THR:N	2.54	0.41
1:C:142:ASP:OD1	1:C:143:SER:N	2.53	0.41
2:F:11:LEU:HD11	2:F:19:LYS:HG3	2.02	0.41
2:F:142:GLU:OE1	2:F:255:ARG:NH1	2.36	0.41
2:H:69:GLU:O	2:H:70:SER:C	2.64	0.41
2:D:83:ASN:OD1	2:D:83:ASN:C	2.63	0.41
1:E:110:THR:O	1:E:110:THR:OG1	2.30	0.41
1:E:138:TYR:HE1	2:F:147:GLU:HA	1.86	0.41
2:F:216:ASP:HA	2:F:217:PRO:HD3	1.95	0.41
1:A:141:ARG:HD3	1:A:141:ARG:HA	1.82	0.41
1:C:156:PHE:CE1	1:C:188:ASN:HB2	2.56	0.41
1:E:67:THR:HG22	1:E:77:SER:HB2	2.03	0.41
1:G:178:LEU:HD13	2:H:184:CYS:HB2	2.03	0.41
2:H:131:LYS:CE	2:H:238:GLN:HB2	2.50	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:131:LYS:NZ	2:B:238:GLN:HG3	2.36	0.41
1:E:142:ASP:OD1	1:E:143:SER:N	2.54	0.41
1:E:157:ASP:OD2	1:E:159:GLN:HG2	2.22	0.41
1:C:209:ILE:O	1:C:209:ILE:HG22	2.20	0.40
2:D:148:ILE:HD12	2:D:148:ILE:H	1.87	0.40
2:F:164:TYR:CD2	2:F:164:TYR:C	2.96	0.40
2:H:137:GLU:O	2:H:160:ALA:HA	2.21	0.40
1:E:154:THR:HG22	1:E:155:ASP:N	2.36	0.40
2:F:183:VAL:HG22	2:F:207:LEU:HD13	2.04	0.40
2:H:143:PRO:HD2	2:H:214:TRP:CE2	2.56	0.40
2:B:195:ALA:HB1	1:C:83:ALA:HB1	2.03	0.40
1:C:165:SER:HB2	1:C:172:ILE:HD11	2.04	0.40
1:G:81:GLN:HE21	1:G:84:ARG:HD2	1.86	0.40
2:F:4:VAL:HG22	2:F:25:GLN:HB3	2.03	0.40
2:F:207:LEU:HD12	2:F:207:LEU:HA	1.81	0.40
1:G:153:PHE:HE2	1:G:162:VAL:HG22	1.86	0.40
2:H:5:THR:CG2	2:H:24:GLU:HB2	2.52	0.40
1:A:56:GLN:CB	1:A:78:LEU:HD23	2.50	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:F:58:TYR:O	2:H:37:ARG:NH2[1_655]	2.10	0.10
2:B:58:TYR:O	2:D:37:ARG:NH2[2_455]	2.16	0.04

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	202/205 (98%)	195 (96%)	7 (4%)	0	100 100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	C	202/205 (98%)	185 (92%)	17 (8%)	0	100	100
1	E	194/205 (95%)	178 (92%)	16 (8%)	0	100	100
1	G	202/205 (98%)	187 (93%)	15 (7%)	0	100	100
2	B	246/249 (99%)	239 (97%)	7 (3%)	0	100	100
2	D	239/249 (96%)	228 (95%)	11 (5%)	0	100	100
2	F	236/249 (95%)	222 (94%)	14 (6%)	0	100	100
2	H	236/249 (95%)	227 (96%)	9 (4%)	0	100	100
All	All	1757/1816 (97%)	1661 (94%)	96 (6%)	0	100	100

There are no Ramachandran outliers to report.

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	181/182 (100%)	177 (98%)	4 (2%)	45	73
1	C	181/182 (100%)	179 (99%)	2 (1%)	65	84
1	E	175/182 (96%)	169 (97%)	6 (3%)	32	60
1	G	181/182 (100%)	174 (96%)	7 (4%)	28	55
2	B	218/219 (100%)	211 (97%)	7 (3%)	34	62
2	D	216/219 (99%)	209 (97%)	7 (3%)	34	62
2	F	215/219 (98%)	209 (97%)	6 (3%)	38	66
2	H	215/219 (98%)	208 (97%)	7 (3%)	33	61
All	All	1582/1604 (99%)	1536 (97%)	46 (3%)	37	65

All (46) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	54	ILE
1	A	92	SER

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Mol	Chain	Res	Type
1	A	165	SER
1	A	173	THR
2	B	21	LEU
2	B	66	SER
2	B	87	LEU
2	B	138	VAL
2	B	190	LEU
2	B	206	ARG
2	B	247	ILE
1	C	164	GLN
1	C	189	SER
2	D	5	THR
2	D	9	LYS
2	D	64	LYS
2	D	113	THR
2	D	128	GLU
2	D	144	SER
2	D	174	VAL
1	E	7	THR
1	E	39	LEU
1	E	53	LEU
1	E	94	SER
1	E	110	THR
1	E	170	VAL
2	F	66	SER
2	F	113	THR
2	F	145	LYS
2	F	147	GLU
2	F	161	THR
2	F	196	LEU
1	G	5	THR
1	G	8	GLN
1	G	13	VAL
1	G	24	THR
1	G	63	GLU
1	G	127	GLN
1	G	173	THR
2	H	22	LYS
2	H	47	LYS
2	H	84	SER
2	H	123	ARG
2	H	190	LEU

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Mol	Chain	Res	Type
2	H	209	VAL
2	H	232	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (32) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	30	GLN
1	A	44	GLN
1	A	164	GLN
1	A	195	ASN
2	B	17	ASN
2	B	44	GLN
2	B	116	GLN
2	B	226	GLN
2	B	233	ASN
2	B	246	GLN
1	C	14	GLN
1	C	30	GLN
1	C	79	ASN
1	C	81	GLN
2	D	10	HIS
2	D	17	ASN
2	D	26	HIS
2	D	92	HIS
2	D	116	GLN
2	D	193	GLN
1	E	2	GLN
1	E	132	ASN
2	F	10	HIS
2	F	83	ASN
2	F	175	ASN
2	F	180	HIS
2	F	219	ASN
2	F	226	GLN
1	G	127	GLN
2	H	26	HIS
2	H	116	GLN
2	H	233	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry ⓘ

14 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	$\# Z > 2$
4	SO4	A	302	-	4,4,4	0.39	0	6,6,6	0.22	0
3	NAG	E	301	1	14,14,15	0.28	0	17,19,21	0.46	0
3	NAG	A	301	1	14,14,15	0.68	0	17,19,21	0.56	0
3	NAG	D	301	2	14,14,15	0.66	1 (7%)	17,19,21	0.72	1 (5%)
3	NAG	B	301	2	14,14,15	0.45	0	17,19,21	0.82	1 (5%)
3	NAG	H	301	2	14,14,15	0.56	0	17,19,21	0.79	1 (5%)
4	SO4	E	303	-	4,4,4	0.35	0	6,6,6	0.37	0
4	SO4	G	302	-	4,4,4	0.38	0	6,6,6	0.27	0
3	NAG	E	302	1	14,14,15	0.70	1 (7%)	17,19,21	0.96	1 (5%)
3	NAG	F	301	2	14,14,15	0.27	0	17,19,21	1.20	1 (5%)
3	NAG	G	301	1	14,14,15	0.50	0	17,19,21	0.52	0
4	SO4	B	302	-	4,4,4	0.27	0	6,6,6	0.24	0
4	SO4	C	301	-	4,4,4	0.44	0	6,6,6	0.27	0
4	SO4	D	302	-	4,4,4	0.29	0	6,6,6	0.28	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	NAG	E	301	1	-	0/6/23/26	0/1/1/1
3	NAG	A	301	1	-	2/6/23/26	0/1/1/1
3	NAG	D	301	2	-	0/6/23/26	0/1/1/1
3	NAG	B	301	2	-	0/6/23/26	0/1/1/1
3	NAG	H	301	2	-	0/6/23/26	0/1/1/1
3	NAG	E	302	1	-	2/6/23/26	0/1/1/1
3	NAG	F	301	2	-	2/6/23/26	0/1/1/1
3	NAG	G	301	1	-	2/6/23/26	0/1/1/1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	302	NAG	C1-C2	2.39	1.55	1.52
3	D	301	NAG	O5-C1	2.18	1.47	1.43

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	F	301	NAG	C1-O5-C5	4.16	117.76	112.19
3	E	302	NAG	C1-O5-C5	3.41	116.76	112.19
3	H	301	NAG	C1-O5-C5	2.59	115.66	112.19
3	D	301	NAG	C1-O5-C5	2.41	115.42	112.19
3	B	301	NAG	C1-O5-C5	2.16	115.09	112.19

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	G	301	NAG	O5-C5-C6-O6
3	E	302	NAG	O5-C5-C6-O6
3	E	302	NAG	C4-C5-C6-O6
3	A	301	NAG	C4-C5-C6-O6
3	G	301	NAG	C4-C5-C6-O6
3	A	301	NAG	O5-C5-C6-O6
3	F	301	NAG	C4-C5-C6-O6
3	F	301	NAG	O5-C5-C6-O6

There are no ring outliers.

3 monomers are involved in 4 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	301	NAG	2	0
3	G	301	NAG	1	0
4	B	302	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	204/205 (99%)	0.22	3 (1%) 72 68	35, 57, 86, 105	0
1	C	204/205 (99%)	0.39	8 (3%) 43 38	38, 65, 102, 111	0
1	E	198/205 (96%)	0.54	12 (6%) 27 23	45, 71, 112, 130	0
1	G	204/205 (99%)	0.71	26 (12%) 8 6	43, 76, 110, 121	0
2	B	248/249 (99%)	0.13	5 (2%) 65 61	32, 55, 92, 113	0
2	D	243/249 (97%)	0.23	5 (2%) 63 59	36, 61, 91, 108	0
2	F	240/249 (96%)	0.65	23 (9%) 13 11	37, 79, 111, 122	0
2	H	240/249 (96%)	0.73	37 (15%) 5 4	39, 77, 117, 146	0
All	All	1781/1816 (98%)	0.45	119 (6%) 24 21	32, 64, 109, 146	0

All (119) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	H	174	VAL	4.2
2	D	154	ALA	4.1
2	H	209	VAL	4.0
2	B	111(A)	ALA	4.0
1	E	137	VAL	4.0
2	H	250	ALA	3.9
2	H	179	VAL	3.7
2	H	157	VAL	3.7
2	H	252	ALA	3.6
2	F	221	PHE	3.6
1	G	137	VAL	3.5
1	G	136	ALA	3.5
2	H	257	ASP	3.5
1	G	152	LEU	3.5
2	H	146	ALA	3.4
2	F	209	VAL	3.4

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Mol	Chain	Res	Type	RSRZ
2	B	111	LEU	3.3
2	F	176	GLY	3.3
2	F	151	THR	3.3
1	G	151	CYS	3.3
1	G	209	ILE	3.2
1	C	205	PHE	3.2
2	F	214	TRP	3.2
1	G	182	SER	3.1
1	E	215	PHE	3.1
2	F	213	PHE	3.1
2	H	156	LEU	3.1
2	H	195	ALA	3.1
2	H	140	VAL	3.1
1	G	192	ALA	3.0
2	H	205	SER	3.0
1	G	130	ILE	3.0
1	G	211	PRO	3.0
1	G	191	VAL	2.9
2	D	256	ALA	2.9
2	H	139	ALA	2.9
2	H	221	PHE	2.9
2	H	172	TRP	2.9
1	E	182	SER	2.9
2	B	112(B)	ILE	2.9
1	C	202	ALA	2.9
1	G	199	PHE	2.8
2	B	112	GLY	2.8
2	F	207	LEU	2.8
1	E	183	MET	2.8
1	E	216	PHE	2.8
2	H	214	TRP	2.7
2	F	139	ALA	2.7
1	C	217	PRO	2.7
1	G	150	VAL	2.7
2	H	151	THR	2.7
2	H	173	TRP	2.7
2	F	208	ARG	2.6
1	A	147	ASP	2.6
1	G	138	TYR	2.6
2	D	110	GLY	2.6
1	G	207	ASN	2.6
2	F	211	ALA	2.6

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Mol	Chain	Res	Type	RSRZ
2	H	154	ALA	2.6
1	E	197	SER	2.5
1	G	183	MET	2.5
1	G	210	ILE	2.5
1	E	178	LEU	2.5
1	C	210	ILE	2.5
2	F	217	PRO	2.5
2	H	182	GLY	2.5
2	H	215	GLN	2.5
1	G	162	VAL	2.5
2	D	2	THR	2.5
2	F	141	PHE	2.4
2	H	236	TRP	2.4
2	F	154	ALA	2.4
2	H	148	ILE	2.4
1	G	140	LEU	2.4
2	F	156	LEU	2.4
1	G	149	PHE	2.4
1	E	142	ASP	2.4
2	H	220	HIS	2.4
2	H	211	ALA	2.4
2	H	256	ALA	2.4
2	F	155	THR	2.4
2	H	167	HIS	2.3
1	G	190	ALA	2.3
1	A	146	SER	2.3
1	E	199	PHE	2.3
1	G	135	PRO	2.3
2	F	158	CYS	2.3
2	F	164	TYR	2.3
2	H	210	SER	2.2
1	G	193	TRP	2.2
1	E	138	TYR	2.2
2	F	159	LEU	2.2
2	H	196	LEU	2.2
2	F	210	SER	2.2
2	H	249	SER	2.2
1	E	193	TRP	2.2
2	B	236	TRP	2.2
1	C	215	PHE	2.2
2	H	143	PRO	2.2
2	H	170	LEU	2.2

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Mol	Chain	Res	Type	RSRZ
1	G	145	SER	2.2
1	G	217	PRO	2.2
1	A	148	LYS	2.2
2	D	209	VAL	2.1
2	F	180	HIS	2.1
1	G	142	ASP	2.1
2	H	181	ASP	2.1
1	E	151	CYS	2.1
2	H	149	SER	2.1
2	H	141	PHE	2.1
1	C	182	SER	2.1
1	C	199	PHE	2.1
2	F	140	VAL	2.1
2	F	172	TRP	2.1
2	H	138	VAL	2.0
2	H	248	VAL	2.0
1	C	10	GLY	2.0
1	G	59	TYR	2.0
2	F	148	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no 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	NAG	E	302	14/15	0.65	0.16	94,104,107,107	0
3	NAG	E	301	14/15	0.67	0.12	84,97,104,105	0
3	NAG	G	301	14/15	0.68	0.11	86,94,104,107	0
3	NAG	A	301	14/15	0.71	0.12	77,81,86,90	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors($\text{\AA}^2$)	Q<0.9
4	SO4	B	302	5/5	0.81	0.14	72,80,95,95	0
4	SO4	D	302	5/5	0.81	0.09	78,80,97,99	0
4	SO4	E	303	5/5	0.84	0.17	69,73,81,86	0
4	SO4	A	302	5/5	0.87	0.15	58,61,65,82	0
4	SO4	G	302	5/5	0.88	0.12	61,64,76,79	0
3	NAG	F	301	14/15	0.92	0.10	31,49,58,67	0
3	NAG	B	301	14/15	0.92	0.09	32,44,52,53	0
4	SO4	C	301	5/5	0.92	0.14	62,63,76,83	0
3	NAG	D	301	14/15	0.94	0.07	31,45,57,61	0
3	NAG	H	301	14/15	0.94	0.06	34,50,53,67	0

6.5 Other polymers [i](#)

There are no such residues in this entry.