



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 18, 2026 – 08:52 AM UTC

PDB ID : 2BJH / pdb_00002bjh
Title : Crystal Structure Of S133A AnFaeA-ferulic acid complex
Authors : Faulds, C.B.; Molina, R.; Gonzalez, R.; Husband, F.; Juge, N.; Sanz-Aparicio, J.; Hermoso, J.A.
Deposited on : 2005-02-02
Resolution : 2.54 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

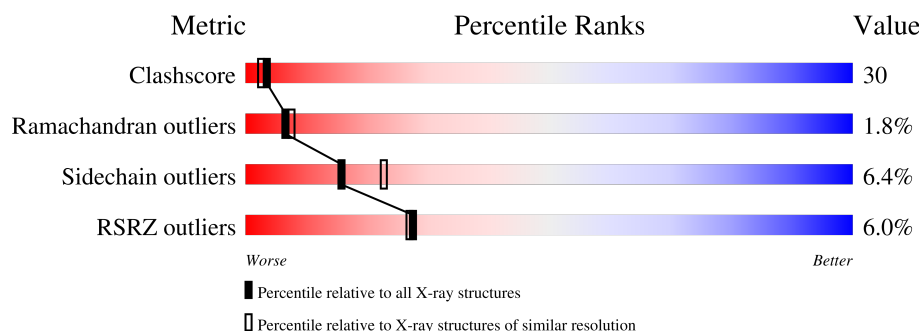
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.54 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))
Clashscore	190562	1120 (2.54-2.54)
Ramachandran outliers	187476	1106 (2.54-2.54)
Sidechain outliers	187428	1106 (2.54-2.54)
RSRZ outliers	180081	1091 (2.54-2.54)

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	260	2% 61% 34% 5%
1	B	260	3% 54% 42% .
1	C	260	13% 47% 46% 7%

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	NAG	C	900	-	-	X	-
4	FER	C	908	-	-	X	-

2 Entry composition [i](#)

There are 5 unique types of molecules in this entry. The entry contains 6478 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.

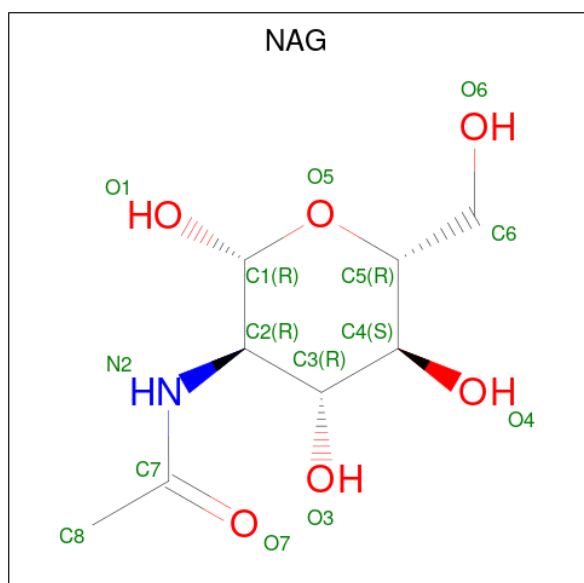
- Molecule 1 is a protein called FERULOYL ESTERASE A.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	260	Total	C	N	O	S	0	0	0
			1991	1239	328	413	11			
1	B	260	Total	C	N	O	S	0	0	0
			1991	1239	328	413	11			
1	C	260	Total	C	N	O	S	0	0	0
			1991	1239	328	413	11			

There are 3 discrepancies between the modelled and reference sequences:

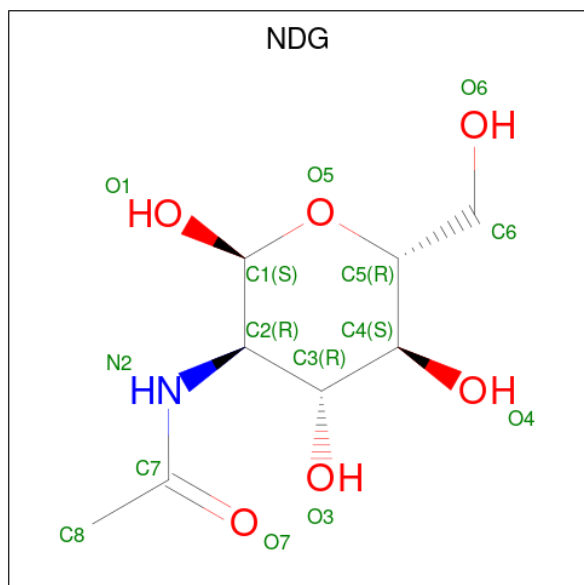
Chain	Residue	Modelled	Actual	Comment	Reference
A	133	ALA	SER	engineered mutation	UNP O42807
B	133	ALA	SER	engineered mutation	UNP O42807
C	133	ALA	SER	engineered mutation	UNP O42807

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



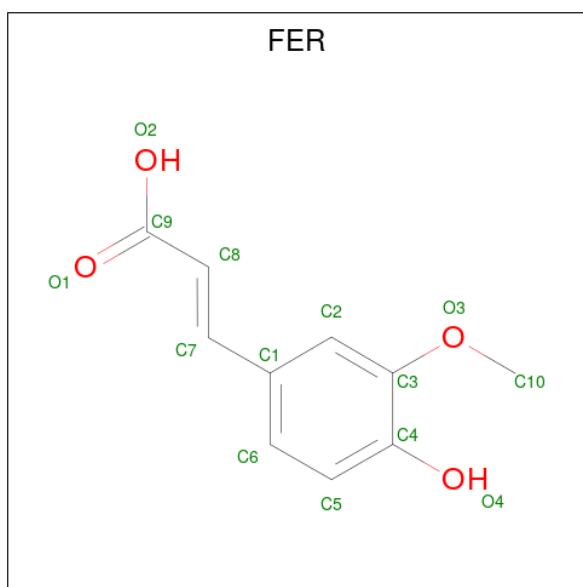
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			15	8	1	6		
2	B	1	Total	C	N	O	0	0
			15	8	1	6		
2	C	1	Total	C	N	O	0	0
			15	8	1	6		

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



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			15	8	1	6		
3	B	1	Total	C	N	O	0	0
			15	8	1	6		
3	C	1	Total	C	N	O	0	0
			15	8	1	6		

- Molecule 4 is 3-(4-HYDROXY-3-METHOXYPHENYL)-2-PROPENOIC ACID (CCD ID: FER) (formula: $C_{10}H_{10}O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			14	10	4		
4	B	1	Total	C	O	0	0
			14	10	4		
4	C	1	Total	C	O	0	0
			14	10	4		

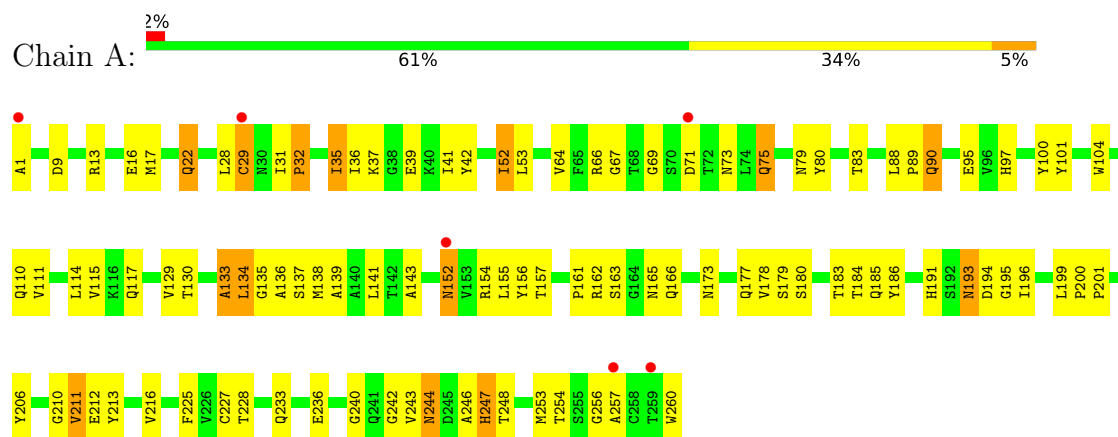
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	157	Total	O	0	0
			157	157		
5	B	139	Total	O	0	0
			139	139		
5	C	77	Total	O	0	0
			77	77		

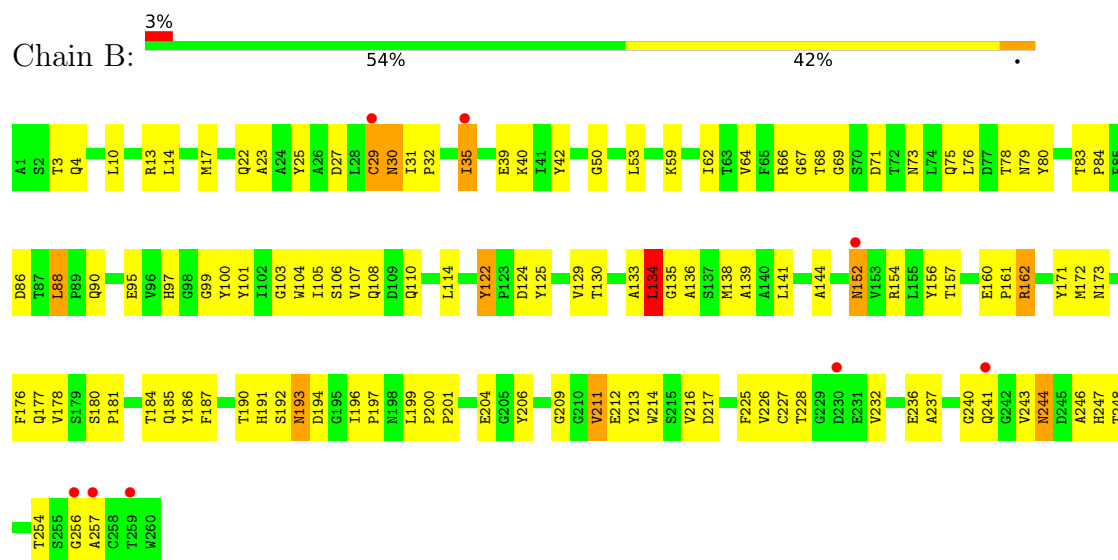
3 Residue-property plots [i](#)

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.

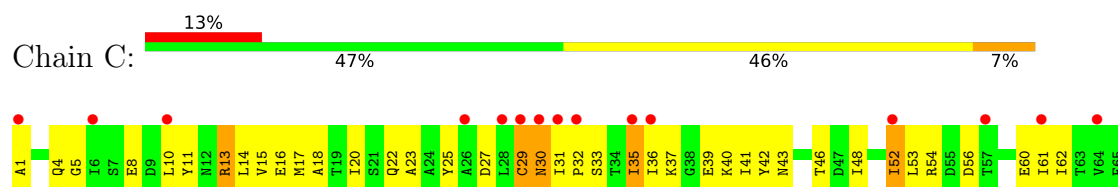
• Molecule 1: FERULOYL ESTERASE A

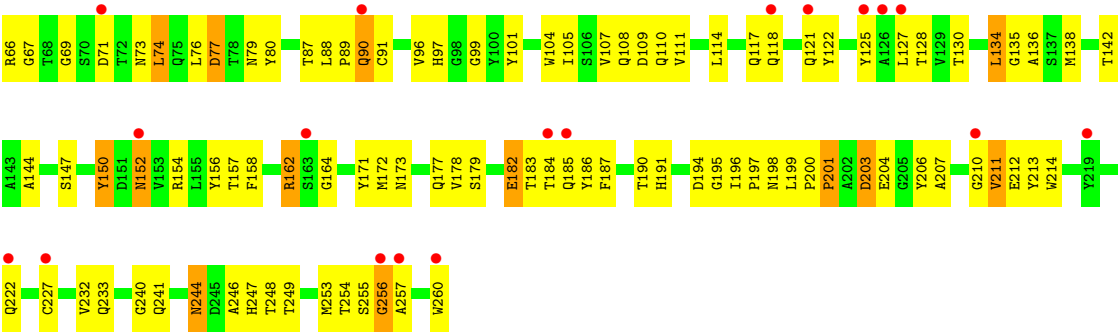


• Molecule 1: FERULOYL ESTERASE A



• Molecule 1: FERULOYL ESTERASE A





4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	46.75Å 130.75Å 76.50Å 90.00° 98.14° 90.00°	Depositor
Resolution (Å)	49.49 – 2.54 49.49 – 2.54	Depositor EDS
% Data completeness (in resolution range)	98.3 (49.49-2.54) 98.3 (49.49-2.54)	Depositor EDS
R_{merge}	0.20	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.10 (at 2.39Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.216 , 0.277 0.211 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å ²)	23.8	Xtriage
Anisotropy	0.733	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 56.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.90	EDS
Total number of atoms	6478	wwPDB-VP
Average B, all atoms (Å ²)	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% 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: NDG, NAG, FER

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	0.44	0/2037	0.95	6/2784 (0.2%)
1	B	0.44	0/2037	0.97	9/2784 (0.3%)
1	C	0.38	0/2037	0.90	4/2784 (0.1%)
All	All	0.42	0/6111	0.94	19/8352 (0.2%)

There are no bond length outliers.

All (19) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	27	ASP	N-CA-C	-7.35	101.73	110.41
1	A	163	SER	N-CA-C	6.00	120.21	113.01
1	A	101	TYR	N-CA-C	-5.94	104.88	111.36
1	A	52	ILE	N-CA-C	-5.83	100.02	108.36
1	C	77	ASP	N-CA-C	5.50	116.96	110.97
1	C	201	PRO	N-CA-C	5.32	118.84	110.80
1	C	162	ARG	N-CA-C	-5.31	100.44	108.67
1	B	192	SER	CB-CA-C	-5.28	110.47	116.54
1	B	122	TYR	CA-C-N	5.24	125.29	119.32
1	B	122	TYR	C-N-CA	5.24	125.29	119.32
1	A	247	HIS	N-CA-C	5.22	116.97	111.28
1	C	240	GLY	N-CA-C	-5.20	105.28	112.25
1	B	88	LEU	CA-C-N	5.17	126.30	119.84
1	B	88	LEU	C-N-CA	5.17	126.30	119.84
1	B	209	GLY	N-CA-C	-5.15	103.67	112.64
1	A	75	GLN	N-CA-C	-5.15	105.67	111.28
1	B	134	LEU	N-CA-C	-5.15	104.74	111.02
1	B	162	ARG	N-CA-C	-5.06	101.86	109.15
1	A	133	ALA	CB-CA-C	-5.03	110.80	116.63

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	1991	0	1836	97	0
1	B	1991	0	1836	114	0
1	C	1991	0	1836	129	0
2	A	15	0	15	4	0
2	B	15	0	15	4	0
2	C	15	0	15	7	0
3	A	15	0	12	2	0
3	B	15	0	12	1	0
3	C	15	0	12	6	0
4	A	14	0	9	0	0
4	B	14	0	9	1	0
4	C	14	0	9	9	0
5	A	157	0	0	6	0
5	B	139	0	0	2	0
5	C	77	0	0	8	0
All	All	6478	0	5616	349	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 30.

All (349) 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:35:ILE:HD11	1:A:53:LEU:HB3	1.30	1.11
1:C:212:GLU:HB3	1:C:227:CYS:HB2	1.46	0.98
1:A:42:TYR:H	1:A:110:GLN:HE22	1.00	0.93
1:A:184:THR:HG22	1:A:186:TYR:H	1.32	0.91
1:B:184:THR:HG22	1:B:186:TYR:H	1.34	0.91
3:C:901:NDG:H4	5:C:2077:HOH:O	1.72	0.89
1:A:42:TYR:N	1:A:110:GLN:HE22	1.71	0.88
1:B:42:TYR:H	1:B:110:GLN:HE22	1.20	0.88
1:B:184:THR:HG21	1:B:211:VAL:HG12	1.53	0.88

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:90:GLN:NE2	1:A:90:GLN:H	1.72	0.87
2:B:900:NAG:O4	3:B:901:NDG:C1	2.24	0.85
1:A:42:TYR:H	1:A:110:GLN:NE2	1.74	0.84
1:A:39:GLU:HG2	1:A:114:LEU:HD22	1.58	0.84
2:A:900:NAG:O4	3:A:901:NDG:C1	2.27	0.83
1:B:152:ASN:HD22	1:B:152:ASN:H	1.23	0.82
1:A:22:GLN:HE21	1:A:22:GLN:HA	1.42	0.82
1:C:96:VAL:HA	1:C:164:GLY:HA2	1.61	0.82
1:A:1:ALA:HB3	1:A:233:GLN:HE22	1.43	0.82
1:B:201:PRO:HG2	1:B:204:GLU:HG3	1.62	0.81
1:C:184:THR:HG22	1:C:186:TYR:H	1.48	0.78
1:C:184:THR:HG21	1:C:211:VAL:HG12	1.65	0.78
1:B:31:ILE:HD12	1:B:35:ILE:HD12	1.65	0.77
1:A:79:ASN:CG	2:A:900:NAG:O1	2.29	0.76
1:B:184:THR:HG21	1:B:211:VAL:CG1	2.16	0.75
1:B:152:ASN:H	1:B:152:ASN:ND2	1.84	0.74
1:A:191:HIS:HD2	1:A:246:ALA:O	1.71	0.73
1:A:212:GLU:HB3	1:A:227:CYS:HB2	1.70	0.73
1:A:89:PRO:HD2	1:A:90:GLN:HE22	1.55	0.72
1:A:90:GLN:H	1:A:90:GLN:HE21	1.36	0.72
1:C:39:GLU:HB3	1:C:52:ILE:HD13	1.72	0.72
1:B:42:TYR:H	1:B:110:GLN:NE2	1.87	0.72
1:A:134:LEU:C	1:A:134:LEU:HD12	2.14	0.71
1:C:88:LEU:HD12	1:C:90:GLN:NE2	2.05	0.71
1:A:193:ASN:HD22	1:A:193:ASN:H	1.40	0.69
1:C:17:MET:HE3	1:C:130:THR:HG21	1.75	0.69
1:A:1:ALA:CB	1:A:233:GLN:HE22	2.05	0.69
1:B:39:GLU:HG2	1:B:114:LEU:HD22	1.74	0.68
1:B:177:GLN:HE21	1:B:185:GLN:CD	2.01	0.68
1:B:152:ASN:HD22	1:B:152:ASN:N	1.84	0.68
1:B:191:HIS:HD2	1:B:246:ALA:O	1.77	0.68
2:C:900:NAG:H61	3:C:901:NDG:HA	1.57	0.68
1:B:103:GLY:O	1:B:106:SER:HB3	1.92	0.68
1:A:179:SER:HB2	5:A:2121:HOH:O	1.94	0.67
1:C:46:THR:OG1	1:C:48:ILE:HG12	1.95	0.67
1:C:8:GLU:HG2	5:C:2008:HOH:O	1.94	0.67
1:A:1:ALA:HB1	1:A:228:THR:O	1.94	0.66
1:B:100:TYR:CE2	1:B:134:LEU:HD22	2.30	0.66
1:C:196:ILE:HG23	4:C:908:FER:H2	1.76	0.66
1:A:184:THR:HG21	1:A:211:VAL:HG13	1.77	0.66
1:A:32:PRO:O	1:A:35:ILE:HG22	1.95	0.65

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:ASN:CG	2:C:900:NAG:O1	2.39	0.65
1:C:46:THR:HB	1:C:76:LEU:HD11	1.78	0.65
1:A:39:GLU:HG2	1:A:114:LEU:CD2	2.26	0.65
1:A:161:PRO:HG3	1:A:196:ILE:CG2	2.27	0.65
1:B:35:ILE:C	1:B:35:ILE:HD13	2.21	0.65
1:C:201:PRO:HG2	1:C:204:GLU:HB2	1.76	0.65
1:B:193:ASN:H	1:B:193:ASN:HD22	1.43	0.65
1:B:35:ILE:HD11	1:B:53:LEU:HD22	1.78	0.64
1:B:248:THR:O	1:B:254:THR:HA	1.98	0.64
1:C:52:ILE:HD12	1:C:52:ILE:N	2.12	0.64
1:B:17:MET:HE3	1:B:130:THR:HG21	1.80	0.63
1:B:133:ALA:HB2	1:B:247:HIS:HE1	1.64	0.63
1:B:193:ASN:O	1:B:240:GLY:HA3	1.98	0.62
1:C:31:ILE:HD12	1:C:35:ILE:HD12	1.80	0.62
1:C:134:LEU:HD12	1:C:135:GLY:N	2.13	0.62
1:C:136:ALA:HB1	1:C:157:THR:HB	1.81	0.62
1:B:88:LEU:HD21	1:B:171:TYR:CZ	2.34	0.62
1:B:40:LYS:HG3	1:B:50:GLY:O	1.99	0.62
1:C:154:ARG:HG3	1:C:154:ARG:HH11	1.65	0.62
1:A:201:PRO:HA	5:A:2132:HOH:O	2.00	0.62
1:C:196:ILE:HD12	4:C:908:FER:H2	1.82	0.61
1:C:196:ILE:HD12	4:C:908:FER:C2	2.30	0.61
1:C:77:ASP:HB3	4:C:908:FER:H5	1.83	0.61
1:B:129:VAL:HG12	1:B:139:ALA:HB1	1.82	0.61
1:C:61:ILE:HD13	1:C:118:GLN:HB3	1.81	0.61
1:A:244:ASN:C	1:A:244:ASN:HD22	2.07	0.60
1:B:104:TRP:HA	1:B:138:MET:HE1	1.82	0.60
1:C:199:LEU:O	4:C:908:FER:H101	2.02	0.60
1:A:104:TRP:HA	1:A:138:MET:HE1	1.84	0.60
1:B:79:ASN:CG	2:B:900:NAG:O1	2.45	0.60
1:C:10:LEU:O	1:C:14:LEU:HG	2.01	0.60
1:B:10:LEU:O	1:B:14:LEU:HG	2.02	0.60
1:C:40:LYS:HE2	1:C:42:TYR:HB2	1.82	0.60
1:B:13:ARG:HD3	1:B:156:TYR:OH	2.02	0.59
1:C:52:ILE:HD11	1:C:114:LEU:HD13	1.82	0.59
1:A:152:ASN:ND2	1:A:152:ASN:H	1.99	0.59
1:C:222:GLN:HB2	5:C:2061:HOH:O	2.02	0.59
1:B:76:LEU:O	1:B:99:GLY:HA3	2.03	0.59
1:C:48:ILE:HG21	1:C:138:MET:HE2	1.83	0.59
1:C:61:ILE:HD13	1:C:118:GLN:CB	2.33	0.59
1:A:9:ASP:HA	5:A:2015:HOH:O	2.02	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:79:ASN:ND2	2:C:900:NAG:O1	2.36	0.58
1:B:244:ASN:H	1:B:244:ASN:ND2	1.99	0.58
1:C:244:ASN:C	1:C:244:ASN:HD22	2.11	0.58
1:C:134:LEU:HD12	1:C:134:LEU:C	2.29	0.58
1:A:22:GLN:HA	1:A:22:GLN:NE2	2.17	0.58
1:B:32:PRO:HB2	1:B:35:ILE:HG22	1.84	0.58
1:B:3:THR:CG2	1:B:226:VAL:HB	2.34	0.57
1:B:3:THR:HG22	1:B:226:VAL:HB	1.86	0.57
1:C:173:ASN:HA	1:C:178:VAL:CG2	2.34	0.57
1:B:181:PRO:HG3	1:B:228:THR:HG21	1.86	0.57
1:B:133:ALA:HB2	1:B:247:HIS:CE1	2.39	0.57
1:C:74:LEU:HA	1:C:77:ASP:HB2	1.85	0.57
1:A:88:LEU:C	1:A:90:GLN:HE21	2.13	0.57
1:B:200:PRO:HG2	1:B:206:TYR:CG	2.39	0.56
1:A:41:ILE:HG23	1:A:110:GLN:HE21	1.70	0.56
1:A:69:GLY:H	1:A:73:ASN:HD22	1.52	0.56
1:A:75:GLN:HB2	5:A:2067:HOH:O	2.05	0.56
1:C:195:GLY:O	1:C:241:GLN:O	2.23	0.56
1:A:194:ASP:OD1	1:A:247:HIS:HB2	2.04	0.56
2:C:900:NAG:H62	3:C:901:NDG:H8C1	1.88	0.56
1:C:69:GLY:H	1:C:73:ASN:HD22	1.54	0.56
1:C:177:GLN:HB2	1:C:183:THR:O	2.05	0.56
1:B:25:TYR:OH	1:B:67:GLY:HA3	2.05	0.56
1:C:23:ALA:HB1	1:C:29:CYS:N	2.21	0.56
1:C:130:THR:HA	1:C:156:TYR:O	2.05	0.55
2:C:900:NAG:C6	3:C:901:NDG:HA	2.19	0.55
1:A:193:ASN:O	1:A:240:GLY:HA3	2.05	0.55
1:B:80:TYR:CD1	1:B:200:PRO:HB3	2.41	0.55
1:A:184:THR:HG22	1:A:186:TYR:N	2.13	0.55
1:C:191:HIS:CD2	1:C:246:ALA:O	2.59	0.55
1:C:35:ILE:HD13	1:C:35:ILE:C	2.32	0.55
1:C:211:VAL:HG13	1:C:213:TYR:HE1	1.72	0.55
1:A:154:ARG:HG3	5:A:2107:HOH:O	2.07	0.55
1:B:211:VAL:HG13	1:B:213:TYR:HE1	1.72	0.54
1:C:35:ILE:HD13	1:C:35:ILE:O	2.07	0.54
1:C:200:PRO:HG2	1:C:206:TYR:CG	2.42	0.54
1:A:89:PRO:HD2	1:A:90:GLN:NE2	2.22	0.54
1:A:200:PRO:HG2	1:A:206:TYR:CG	2.43	0.54
1:C:191:HIS:HD2	1:C:246:ALA:O	1.91	0.54
1:C:62:ILE:HA	1:C:128:THR:O	2.07	0.54
2:C:900:NAG:O4	3:C:901:NDG:C7	2.55	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:144:ALA:HB2	1:B:172:MET:SD	2.48	0.54
1:B:157:THR:HB	1:B:160:GLU:HB2	1.89	0.54
1:B:161:PRO:HG3	1:B:196:ILE:CG2	2.38	0.54
1:A:133:ALA:HB2	1:A:247:HIS:HE1	1.73	0.53
1:B:42:TYR:N	1:B:110:GLN:HE22	1.99	0.53
1:C:52:ILE:CD1	1:C:114:LEU:HD13	2.36	0.53
1:B:68:THR:HA	1:B:73:ASN:HD22	1.73	0.53
1:B:69:GLY:H	1:B:73:ASN:HD22	1.55	0.53
1:C:248:THR:O	1:C:254:THR:HA	2.08	0.53
1:C:10:LEU:HD11	1:C:187:PHE:CE2	2.43	0.53
1:A:28:LEU:HD12	1:A:37:LYS:HE3	1.90	0.53
1:B:29:CYS:O	1:B:30:ASN:C	2.51	0.53
1:C:90:GLN:NE2	1:C:90:GLN:H	2.07	0.53
1:B:67:GLY:HA2	1:B:135:GLY:H	1.73	0.53
1:B:4:GLN:HB2	1:B:225:PHE:CE2	2.44	0.53
1:C:196:ILE:HG23	4:C:908:FER:H103	1.91	0.53
1:C:31:ILE:HB	1:C:32:PRO:HD2	1.90	0.53
1:A:133:ALA:HB2	1:A:247:HIS:CE1	2.44	0.52
1:C:25:TYR:HA	1:C:66:ARG:HG3	1.90	0.52
1:B:173:ASN:HB3	1:B:178:VAL:HB	1.90	0.52
1:A:166:GLN:HB2	5:A:2112:HOH:O	2.08	0.52
1:C:201:PRO:HB2	1:C:203:ASP:OD1	2.09	0.52
1:A:1:ALA:HB3	1:A:233:GLN:NE2	2.20	0.52
1:B:84:PRO:HB2	1:B:86:ASP:OD1	2.09	0.52
1:C:36:ILE:N	1:C:36:ILE:HD12	2.24	0.52
1:C:107:VAL:O	1:C:111:VAL:HG23	2.09	0.52
1:A:95:GLU:HB2	1:A:165:ASN:ND2	2.25	0.52
1:A:111:VAL:O	1:A:115:VAL:HG23	2.09	0.52
1:C:41:ILE:HA	1:C:110:GLN:OE1	2.10	0.52
1:A:41:ILE:HD13	1:A:111:VAL:HG22	1.92	0.52
1:C:37:LYS:HA	1:C:53:LEU:HD22	1.91	0.52
1:C:88:LEU:HD21	1:C:171:TYR:CZ	2.44	0.52
1:A:177:GLN:HE21	1:A:185:GLN:NE2	2.08	0.51
1:C:11:TYR:O	1:C:15:VAL:HG23	2.11	0.51
1:A:90:GLN:HE21	1:A:90:GLN:N	2.06	0.51
1:A:136:ALA:HB1	1:A:157:THR:HB	1.92	0.51
1:C:29:CYS:O	1:C:31:ILE:N	2.44	0.51
1:C:35:ILE:HD11	1:C:53:LEU:HD13	1.92	0.51
1:C:178:VAL:HG21	1:C:210:GLY:HA2	1.92	0.51
1:B:154:ARG:CD	1:B:185:GLN:HB3	2.40	0.51
1:C:196:ILE:HG23	4:C:908:FER:C10	2.41	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:196:ILE:HD11	1:B:243:VAL:CG2	2.41	0.51
1:A:178:VAL:HG13	1:A:211:VAL:HG12	1.93	0.51
1:C:90:GLN:H	1:C:90:GLN:CD	2.19	0.51
1:B:79:ASN:CG	2:B:900:NAG:C1	2.85	0.50
1:C:40:LYS:HG3	5:C:2025:HOH:O	2.10	0.50
1:A:129:VAL:HG12	1:A:139:ALA:HB1	1.93	0.50
1:B:177:GLN:HE21	1:B:185:GLN:NE2	2.10	0.50
1:C:87:THR:C	1:C:89:PRO:HD3	2.35	0.50
1:B:31:ILE:HB	1:B:32:PRO:HD2	1.94	0.50
1:B:191:HIS:HE1	1:B:217:ASP:OD2	1.95	0.50
1:A:243:VAL:HG13	1:A:247:HIS:HD2	1.77	0.50
1:B:173:ASN:HA	1:B:178:VAL:CG2	2.41	0.50
1:C:178:VAL:CG2	1:C:210:GLY:HA2	2.42	0.50
1:C:191:HIS:CD2	1:C:249:THR:O	2.65	0.50
1:A:35:ILE:HD13	1:A:35:ILE:C	2.37	0.50
1:A:83:THR:HB	2:A:900:NAG:H62	1.94	0.50
1:C:40:LYS:HE3	5:C:2025:HOH:O	2.11	0.50
1:A:137:SER:OG	1:A:161:PRO:HD2	2.12	0.49
1:B:212:GLU:HB3	1:B:227:CYS:HB2	1.94	0.49
1:C:184:THR:HG21	1:C:211:VAL:CG1	2.38	0.49
1:B:83:THR:O	1:B:95:GLU:HA	2.13	0.49
1:C:88:LEU:HD12	1:C:90:GLN:HE21	1.74	0.49
1:B:193:ASN:HD22	1:B:193:ASN:N	2.04	0.49
1:B:196:ILE:HD11	1:B:243:VAL:HG23	1.94	0.48
1:C:162:ARG:NH1	1:C:198:ASN:O	2.44	0.48
1:A:134:LEU:C	1:A:134:LEU:CD1	2.86	0.48
2:A:900:NAG:H61	3:A:901:NDG:O1	2.14	0.48
1:B:79:ASN:ND2	2:B:900:NAG:H1	2.29	0.48
1:C:80:TYR:CD1	1:C:200:PRO:HB3	2.48	0.48
1:A:13:ARG:NH1	1:A:156:TYR:OH	2.46	0.48
1:B:196:ILE:N	1:B:197:PRO:HD2	2.29	0.48
1:C:16:GLU:O	1:C:20:ILE:HG13	2.12	0.48
1:C:150:TYR:CD1	1:C:150:TYR:N	2.80	0.48
1:B:193:ASN:H	1:B:193:ASN:ND2	2.11	0.48
1:A:134:LEU:HD12	1:A:134:LEU:O	2.13	0.48
1:B:211:VAL:HG13	1:B:213:TYR:CE1	2.47	0.48
1:C:48:ILE:CG2	1:C:138:MET:HE2	2.43	0.48
1:C:194:ASP:OD1	1:C:197:PRO:HD2	2.13	0.48
1:A:211:VAL:HG22	1:A:213:TYR:CE1	2.48	0.48
1:C:195:GLY:C	1:C:241:GLN:O	2.57	0.48
1:C:61:ILE:HB	1:C:127:LEU:HD23	1.96	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:97:HIS:CG	1:A:100:TYR:CD1	3.01	0.47
1:B:184:THR:HG22	1:B:185:GLN:N	2.30	0.47
1:C:122:TYR:HB3	1:C:125:TYR:CG	2.49	0.47
1:B:17:MET:HE3	1:B:130:THR:CG2	2.44	0.47
1:B:100:TYR:OH	4:B:908:FER:H103	2.15	0.47
1:B:66:ARG:HH21	1:B:73:ASN:HD21	1.62	0.47
1:C:257:ALA:HA	5:C:2073:HOH:O	2.14	0.47
1:B:39:GLU:HG2	1:B:114:LEU:CD2	2.42	0.47
1:B:104:TRP:C	1:B:106:SER:H	2.23	0.47
1:B:173:ASN:HA	1:B:178:VAL:HG23	1.97	0.47
1:C:88:LEU:HD21	1:C:171:TYR:CE2	2.49	0.47
1:B:216:VAL:HG12	1:B:217:ASP:N	2.28	0.47
1:B:244:ASN:H	1:B:244:ASN:HD22	1.62	0.47
1:C:74:LEU:O	1:C:77:ASP:HB2	2.15	0.47
1:A:17:MET:HE3	1:A:130:THR:HG21	1.96	0.47
1:A:67:GLY:HA2	1:A:135:GLY:H	1.79	0.47
1:B:243:VAL:HG12	1:B:243:VAL:O	2.15	0.47
1:A:180:SER:HB3	1:A:183:THR:OG1	2.14	0.46
1:C:18:ALA:HB2	1:C:158:PHE:CE2	2.50	0.46
1:C:104:TRP:CZ3	1:C:142:THR:HA	2.51	0.46
1:C:4:GLN:HG3	1:C:5:GLY:N	2.31	0.46
1:B:31:ILE:CD1	1:B:35:ILE:HD12	2.41	0.46
1:B:67:GLY:HA2	1:B:134:LEU:HB3	1.97	0.46
1:B:133:ALA:O	1:B:134:LEU:C	2.58	0.46
1:C:89:PRO:C	1:C:91:CYS:N	2.70	0.46
1:C:241:GLN:NE2	5:C:2067:HOH:O	2.48	0.46
1:C:42:TYR:CG	1:C:43:ASN:N	2.83	0.46
1:A:31:ILE:O	1:A:32:PRO:C	2.58	0.46
1:A:143:ALA:HB2	1:A:155:LEU:HD13	1.98	0.45
1:A:152:ASN:N	1:A:152:ASN:HD22	2.13	0.45
1:C:232:VAL:O	1:C:233:GLN:HB3	2.15	0.45
1:B:176:PHE:O	1:B:185:GLN:HG3	2.16	0.45
1:B:23:ALA:HB1	1:B:29:CYS:N	2.31	0.45
1:C:1:ALA:HB3	1:C:233:GLN:HE22	1.81	0.45
1:A:35:ILE:CD1	1:A:53:LEU:HB3	2.22	0.45
1:A:52:ILE:HD12	1:A:52:ILE:N	2.31	0.45
1:A:28:LEU:HD12	1:A:37:LYS:HG3	1.99	0.45
1:C:54:ARG:HE	1:C:56:ASP:CG	2.23	0.45
1:A:42:TYR:N	1:A:110:GLN:NE2	2.48	0.45
1:B:129:VAL:HG12	1:B:139:ALA:CB	2.46	0.45
1:C:35:ILE:CD1	1:C:53:LEU:HD13	2.47	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:36:ILE:N	1:A:36:ILE:HD12	2.32	0.45
1:B:154:ARG:HD3	1:B:185:GLN:HB3	1.98	0.45
1:A:133:ALA:O	1:A:136:ALA:N	2.49	0.45
1:C:108:GLN:O	1:C:109:ASP:C	2.60	0.45
1:B:200:PRO:HG2	1:B:206:TYR:CD1	2.52	0.45
1:A:66:ARG:HG2	1:A:67:GLY:N	2.32	0.44
1:A:184:THR:OG1	1:A:211:VAL:HG11	2.17	0.44
1:B:154:ARG:HD2	1:B:185:GLN:HB3	1.99	0.44
1:B:162:ARG:HH21	1:B:236:GLU:CD	2.25	0.44
1:C:134:LEU:C	1:C:134:LEU:CD1	2.90	0.44
1:B:101:TYR:O	1:B:105:ILE:HG23	2.18	0.44
1:C:117:GLN:HG2	1:C:121:GLN:HE21	1.82	0.44
1:A:161:PRO:HG3	1:A:196:ILE:HG22	2.00	0.44
1:B:104:TRP:C	1:B:106:SER:N	2.73	0.44
1:B:194:ASP:OD1	1:B:247:HIS:HB2	2.17	0.44
1:A:152:ASN:H	1:A:152:ASN:HD22	1.66	0.44
1:C:173:ASN:HA	1:C:178:VAL:HG23	2.00	0.44
1:B:244:ASN:HB2	5:B:2128:HOH:O	2.17	0.44
1:C:61:ILE:HD12	1:C:127:LEU:HD21	2.00	0.44
1:C:66:ARG:HG2	1:C:67:GLY:N	2.32	0.44
1:A:97:HIS:HB3	1:A:100:TYR:HB2	1.99	0.43
1:C:101:TYR:O	1:C:105:ILE:HG12	2.18	0.43
1:A:195:GLY:O	1:A:199:LEU:HD23	2.17	0.43
1:B:88:LEU:HG	1:B:90:GLN:HG2	1.99	0.43
1:A:152:ASN:ND2	1:A:152:ASN:N	2.61	0.43
1:C:77:ASP:CB	4:C:908:FER:H5	2.48	0.43
1:C:154:ARG:HG3	1:C:154:ARG:NH1	2.31	0.43
1:A:162:ARG:HH21	1:A:236:GLU:CD	2.27	0.43
1:A:178:VAL:HG21	1:A:210:GLY:HA2	2.00	0.43
1:C:194:ASP:OD2	1:C:247:HIS:HA	2.18	0.43
1:C:254:THR:HG22	1:C:255:SER:N	2.34	0.43
1:C:162:ARG:HD3	1:C:207:ALA:HA	2.01	0.43
1:B:196:ILE:HG13	1:B:247:HIS:CE1	2.54	0.43
1:C:22:GLN:HA	1:C:22:GLN:NE2	2.34	0.43
1:C:89:PRO:C	1:C:91:CYS:H	2.27	0.43
1:B:244:ASN:HD22	1:B:244:ASN:N	2.17	0.43
1:C:253:MET:HB3	1:C:260:TRP:CE3	2.53	0.43
1:A:253:MET:HB3	1:A:260:TRP:CZ3	2.54	0.43
1:B:31:ILE:HB	1:B:32:PRO:CD	2.48	0.43
1:C:52:ILE:HD12	1:C:52:ILE:H	1.84	0.43
1:C:194:ASP:HA	1:C:244:ASN:HD21	1.83	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:41:ILE:HA	1:A:110:GLN:NE2	2.34	0.42
1:B:71:ASP:O	1:B:75:GLN:HG3	2.19	0.42
1:B:244:ASN:HD22	1:B:244:ASN:C	2.27	0.42
1:C:144:ALA:O	1:C:147:SER:HB2	2.19	0.42
1:C:178:VAL:O	1:C:179:SER:C	2.61	0.42
1:C:184:THR:HG22	1:C:185:GLN:N	2.32	0.42
1:B:187:PHE:CD1	1:B:187:PHE:N	2.87	0.42
1:B:100:TYR:HE2	1:B:134:LEU:HD22	1.80	0.42
1:C:13:ARG:HH22	1:C:60:GLU:CD	2.27	0.42
1:C:190:THR:OG1	1:C:214:TRP:HA	2.19	0.42
1:C:196:ILE:HG13	1:C:247:HIS:CE1	2.54	0.42
1:A:22:GLN:HE21	1:A:22:GLN:CA	2.23	0.42
1:A:35:ILE:HD13	1:A:36:ILE:N	2.34	0.42
1:B:53:LEU:HB2	1:B:62:ILE:HB	2.02	0.42
1:C:104:TRP:HA	1:C:138:MET:HE1	2.02	0.42
1:C:79:ASN:O	1:C:97:HIS:CE1	2.73	0.42
1:C:152:ASN:C	1:C:152:ASN:HD22	2.28	0.42
1:B:133:ALA:O	1:B:136:ALA:HB3	2.19	0.42
1:C:255:SER:O	1:C:256:GLY:C	2.62	0.42
1:A:253:MET:HB3	1:A:260:TRP:CE3	2.54	0.41
1:B:162:ARG:HB3	1:B:206:TYR:HB3	2.01	0.41
1:A:89:PRO:CD	1:A:90:GLN:NE2	2.83	0.41
1:A:244:ASN:C	1:A:244:ASN:ND2	2.76	0.41
1:A:17:MET:CE	1:A:130:THR:HB	2.51	0.41
1:B:122:TYR:HB3	1:B:125:TYR:CD1	2.56	0.41
1:B:178:VAL:O	1:B:178:VAL:HG12	2.20	0.41
1:B:180:SER:HA	1:B:181:PRO:HD3	1.92	0.41
1:C:77:ASP:CG	4:C:908:FER:H6	2.45	0.41
1:B:78:THR:HG21	5:B:2059:HOH:O	2.19	0.41
1:B:17:MET:HG3	1:B:156:TYR:CG	2.56	0.41
1:B:83:THR:HA	1:B:84:PRO:HD3	1.95	0.41
1:C:173:ASN:HB3	1:C:178:VAL:HB	2.03	0.41
1:A:248:THR:O	1:A:254:THR:HA	2.20	0.41
1:B:97:HIS:CG	1:B:100:TYR:CD1	3.08	0.41
1:B:190:THR:OG1	1:B:214:TRP:HA	2.21	0.41
1:C:41:ILE:HD13	1:C:111:VAL:HG22	2.02	0.41
1:A:225:PHE:CD1	1:A:225:PHE:N	2.88	0.41
1:B:232:VAL:HG12	1:B:237:ALA:HB2	2.02	0.41
1:B:22:GLN:HA	1:B:22:GLN:HE21	1.86	0.40
1:C:144:ALA:HB2	1:C:172:MET:SD	2.60	0.40
1:C:182:GLU:HG2	5:C:2053:HOH:O	2.20	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:88:LEU:C	1:A:90:GLN:NE2	2.79	0.40
1:A:243:VAL:HG13	1:A:247:HIS:CD2	2.55	0.40
1:C:177:GLN:C	1:C:179:SER:N	2.79	0.40
1:B:107:VAL:O	1:B:108:GLN:C	2.64	0.40
2:C:900:NAG:O4	3:C:901:NDG:N2	2.54	0.40
1:A:80:TYR:CD1	1:A:200:PRO:HB3	2.56	0.40
1:A:193:ASN:H	1:A:193:ASN:ND2	2.15	0.40
1:B:110:GLN:O	1:B:114:LEU:HG	2.21	0.40
1:C:97:HIS:ND1	1:C:99:GLY:N	2.68	0.40
1:A:173:ASN:OD1	1:A:178:VAL:HG11	2.22	0.40
1:B:66:ARG:HH21	1:B:73:ASN:ND2	2.19	0.40
1:C:1:ALA:CB	1:C:233:GLN:HE22	2.35	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	258/260 (99%)	239 (93%)	14 (5%)	5 (2%)	6	7
1	B	258/260 (99%)	235 (91%)	19 (7%)	4 (2%)	7	9
1	C	258/260 (99%)	227 (88%)	26 (10%)	5 (2%)	6	7
All	All	774/780 (99%)	701 (91%)	59 (8%)	14 (2%)	6	7

All (14) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	29	CYS
1	B	241	GLN
1	B	257	ALA
1	C	29	CYS
1	C	30	ASN

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Mol	Chain	Res	Type
1	A	256	GLY
1	B	30	ASN
1	C	256	GLY
1	A	257	ALA
1	B	256	GLY
1	A	32	PRO
1	C	33	SER
1	C	182	GLU
1	A	242	GLY

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	214/214 (100%)	199 (93%)	15 (7%)	14	19
1	B	214/214 (100%)	202 (94%)	12 (6%)	19	29
1	C	214/214 (100%)	200 (94%)	14 (6%)	15	22
All	All	642/642 (100%)	601 (94%)	41 (6%)	16	23

All (41) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	16	GLU
1	A	22	GLN
1	A	29	CYS
1	A	35	ILE
1	A	64	VAL
1	A	71	ASP
1	A	90	GLN
1	A	117	GLN
1	A	134	LEU
1	A	141	LEU
1	A	152	ASN
1	A	193	ASN
1	A	211	VAL

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Mol	Chain	Res	Type
1	A	216	VAL
1	A	244	ASN
1	B	29	CYS
1	B	35	ILE
1	B	59	LYS
1	B	64	VAL
1	B	124	ASP
1	B	134	LEU
1	B	141	LEU
1	B	152	ASN
1	B	193	ASN
1	B	199	LEU
1	B	211	VAL
1	B	244	ASN
1	C	13	ARG
1	C	27	ASP
1	C	30	ASN
1	C	35	ILE
1	C	52	ILE
1	C	71	ASP
1	C	74	LEU
1	C	90	GLN
1	C	134	LEU
1	C	150	TYR
1	C	152	ASN
1	C	203	ASP
1	C	211	VAL
1	C	244	ASN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (45) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	ASN
1	A	22	GLN
1	A	30	ASN
1	A	73	ASN
1	A	90	GLN
1	A	108	GLN
1	A	110	GLN
1	A	117	GLN
1	A	152	ASN
1	A	177	GLN

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Mol	Chain	Res	Type
1	A	191	HIS
1	A	193	ASN
1	A	233	GLN
1	A	244	ASN
1	A	247	HIS
1	B	12	ASN
1	B	22	GLN
1	B	73	ASN
1	B	90	GLN
1	B	92	ASN
1	B	108	GLN
1	B	110	GLN
1	B	152	ASN
1	B	177	GLN
1	B	185	GLN
1	B	191	HIS
1	B	193	ASN
1	B	198	ASN
1	B	233	GLN
1	B	244	ASN
1	B	247	HIS
1	C	4	GLN
1	C	30	ASN
1	C	73	ASN
1	C	79	ASN
1	C	90	GLN
1	C	92	ASN
1	C	110	GLN
1	C	121	GLN
1	C	152	ASN
1	C	185	GLN
1	C	191	HIS
1	C	193	ASN
1	C	233	GLN
1	C	244	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

9 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	NAG	A	900	-	15,15,15	0.47	0	21,21,21	0.76	0
3	NDG	B	901	-	15,15,15	0.40	0	21,21,21	0.69	0
4	FER	A	908	-	14,14,14	1.64	3 (21%)	18,18,18	0.70	0
2	NAG	B	900	-	15,15,15	0.45	0	21,21,21	0.84	0
4	FER	B	908	-	14,14,14	1.51	2 (14%)	18,18,18	0.67	0
3	NDG	A	901	-	15,15,15	0.37	0	21,21,21	0.65	0
4	FER	C	908	-	14,14,14	1.37	3 (21%)	18,18,18	0.59	0
2	NAG	C	900	-	15,15,15	0.37	0	21,21,21	0.63	0
3	NDG	C	901	-	15,15,15	0.45	0	21,21,21	0.56	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	NAG	A	900	-	-	2/6/26/26	0/1/1/1
3	NDG	B	901	-	-	6/6/26/26	0/1/1/1
4	FER	A	908	-	-	0/7/7/7	0/1/1/1
2	NAG	B	900	-	-	3/6/26/26	0/1/1/1
4	FER	B	908	-	-	0/7/7/7	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	NDG	A	901	-	-	2/6/26/26	0/1/1/1
4	FER	C	908	-	-	0/7/7/7	0/1/1/1
2	NAG	C	900	-	-	4/6/26/26	0/1/1/1
3	NDG	C	901	-	-	2/6/26/26	0/1/1/1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	908	FER	C2-C3	3.28	1.44	1.38
4	A	908	FER	C2-C3	3.18	1.44	1.38
4	B	908	FER	C2-C3	3.14	1.44	1.38
4	A	908	FER	C6-C1	2.28	1.44	1.39
4	C	908	FER	O2-C9	-2.26	1.24	1.30
4	A	908	FER	C6-C5	2.17	1.42	1.38
4	B	908	FER	C6-C1	2.04	1.43	1.39
4	C	908	FER	C6-C5	2.02	1.42	1.38

There are no bond angle outliers.

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	C	900	NAG	C8-C7-N2-C2
2	C	900	NAG	O7-C7-N2-C2
2	C	900	NAG	O5-C5-C6-O6
3	A	901	NDG	O5-C5-C6-O6
3	A	901	NDG	C4-C5-C6-O6
2	C	900	NAG	C4-C5-C6-O6
3	B	901	NDG	O5-C5-C6-O6
3	C	901	NDG	O5-C5-C6-O6
3	B	901	NDG	C4-C5-C6-O6
2	B	900	NAG	C8-C7-N2-C2
3	B	901	NDG	C8-C7-N2-C2
2	B	900	NAG	O7-C7-N2-C2
3	B	901	NDG	O7-C7-N2-C2
3	C	901	NDG	C4-C5-C6-O6
3	B	901	NDG	C3-C2-N2-C7
2	A	900	NAG	O5-C5-C6-O6
2	A	900	NAG	C4-C5-C6-O6
2	B	900	NAG	C3-C2-N2-C7

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Mol	Chain	Res	Type	Atoms
3	B	901	NDG	C1-C2-N2-C7

There are no ring outliers.

8 monomers are involved in 26 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	900	NAG	4	0
3	B	901	NDG	1	0
2	B	900	NAG	4	0
4	B	908	FER	1	0
3	A	901	NDG	2	0
4	C	908	FER	9	0
2	C	900	NAG	7	0
3	C	901	NDG	6	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	260/260 (100%)	-0.05	6 (2%) 61 62	6, 19, 36, 45	0
1	B	260/260 (100%)	0.03	8 (3%) 51 52	7, 20, 38, 52	0
1	C	260/260 (100%)	1.07	33 (12%) 8 7	21, 44, 63, 68	0
All	All	780/780 (100%)	0.35	47 (6%) 27 27	6, 24, 57, 68	0

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	257	ALA	5.3
1	C	152	ASN	4.6
1	C	31	ILE	4.3
1	C	210	GLY	3.8
1	C	260	TRP	3.6
1	C	256	GLY	3.4
1	B	152	ASN	3.3
1	B	256	GLY	3.3
1	B	230	ASP	3.2
1	C	30	ASN	3.0
1	C	163	SER	2.9
1	C	121	GLN	2.9
1	B	257	ALA	2.8
1	C	29	CYS	2.8
1	C	127	LEU	2.8
1	B	241	GLN	2.8
1	C	61	ILE	2.7
1	C	125	TYR	2.7
1	A	29	CYS	2.7
1	C	52	ILE	2.7
1	C	64	VAL	2.7
1	C	26	ALA	2.7
1	A	152	ASN	2.6

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Mol	Chain	Res	Type	RSRZ
1	C	28	LEU	2.5
1	A	257	ALA	2.5
1	C	71	ASP	2.5
1	C	118	GLN	2.5
1	B	259	THR	2.5
1	A	71	ASP	2.4
1	C	6	ILE	2.4
1	C	1	ALA	2.4
1	C	10	LEU	2.3
1	C	185	GLN	2.3
1	A	259	THR	2.3
1	C	184	THR	2.3
1	C	57	THR	2.2
1	C	35	ILE	2.2
1	C	32	PRO	2.2
1	B	35	ILE	2.2
1	C	126	ALA	2.1
1	A	1	ALA	2.1
1	C	36	ILE	2.1
1	C	227	CYS	2.1
1	C	90	GLN	2.1
1	C	222	GLN	2.1
1	C	219	TYR	2.0
1	B	29	CYS	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	NDG	C	901	15/15	0.60	0.27	85,90,91,91	0
4	FER	A	908	14/14	0.80	0.25	76,78,83,83	0
3	NDG	B	901	15/15	0.82	0.14	24,29,31,38	0
4	FER	C	908	14/14	0.82	0.28	78,81,82,82	0
4	FER	B	908	14/14	0.85	0.22	71,73,77,78	0
3	NDG	A	901	15/15	0.85	0.16	33,36,41,42	0
2	NAG	C	900	15/15	0.89	0.12	38,40,42,43	0
2	NAG	B	900	15/15	0.91	0.11	25,27,30,30	0
2	NAG	A	900	15/15	0.93	0.10	17,20,24,26	0

6.5 Other polymers [i](#)

There are no such residues in this entry.