



Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 09:57 PM GMT

PDB ID : 1RE3
Title : Crystal Structure of Fragment D of BbetaD398A Fibrinogen with the Peptide
Ligand Gly-His-Arg-Pro-Amide
Authors : Kostelansky, M.S.; Betts, L.; Gorkun, O.V.; Lord, S.T.
Deposited on : 2003-11-06
Resolution : 2.45 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
<http://wwpdb.org/validation/2016/XrayValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.7 (RC4), CSD as536be (2015)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20026688
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)
Refmac : 5.8.0135
CCP4 : 6.5.0
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : trunk26865

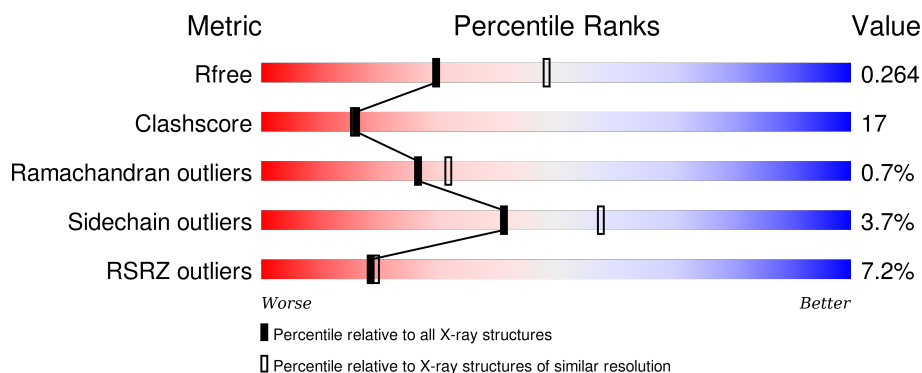
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.45 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.





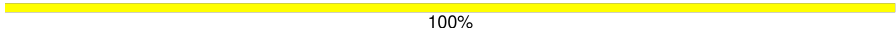
Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	4776 (2.50-2.42)
Clashscore	102246	1030 (2.48-2.44)
Ramachandran outliers	100387	1024 (2.48-2.44)
Sidechain outliers	100360	1024 (2.48-2.44)
RSRZ outliers	91569	4787 (2.50-2.42)

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 on the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria. 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	66	<div> <div>23%</div> <div>41%</div> <div>48%</div> <div>9%</div> </div>
1	D	66	<div> <div>9%</div> <div>53%</div> <div>36%</div> <div>6%</div> <div>5%</div> </div>
2	B	313	<div> <div>8%</div> <div>67%</div> <div>25%</div> <div>6%</div> </div>
2	E	313	<div> <div>6%</div> <div>66%</div> <div>27%</div> <div>•</div> </div>
3	C	311	<div> <div>5%</div> <div>73%</div> <div>18%</div> <div>7%</div> </div>

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Mol	Chain	Length	Quality of chain
3	F	311	
4	G	4	
4	H	4	

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 10794 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 Fibrinogen alpha/alpha-E chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	60	Total	C	N	O	S	0	0	0
			490	300	94	93	3			
1	D	63	Total	C	N	O	S	0	0	0
			516	317	99	97	3			

- Molecule 2 is a protein called Fibrinogen beta chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	294	Total	C	N	O	S	0	0	0
			2359	1474	417	446	22			
2	E	301	Total	C	N	O	S	0	0	0
			2410	1505	426	457	22			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	398	ALA	ASP	ENGINEERED	UNP P02675
E	398	ALA	ASP	ENGINEERED	UNP P02675

- Molecule 3 is a protein called Fibrinogen gamma chain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	288	Total	C	N	O	S	0	0	0
			2312	1468	390	443	11			
3	F	291	Total	C	N	O	S	0	0	0
			2336	1481	395	449	11			

- Molecule 4 is a protein called GHRP peptide.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	G	4	Total	C	N	O	0	0	0
			33	19	9	5			

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace
4	H	4	Total	C	N	O	0	0	0
			33	19	9	5			

- Molecule 5 is a polymer of unknown type called SUGAR (NAG-NAG).

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	B	2	Total	C	N	O	0	0
			28	16	2	10		
5	E	2	Total	C	N	O	0	0
			28	16	2	10		

- Molecule 6 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	B	1	Total	Ca	0	0
			1	1		
6	C	2	Total	Ca	0	0
			2	2		
6	F	1	Total	Ca	0	0
			1	1		
6	E	1	Total	Ca	0	0
			1	1		

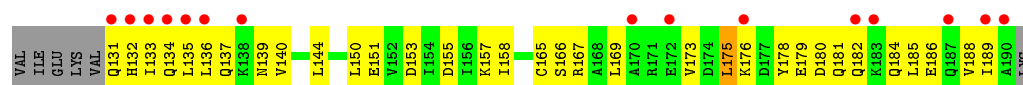
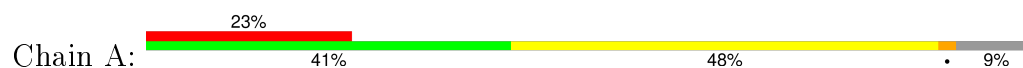
- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
7	A	2	Total	O	0	0
			2	2		
7	B	63	Total	O	0	0
			63	63		
7	C	54	Total	O	0	0
			54	54		
7	D	9	Total	O	0	0
			9	9		
7	E	63	Total	O	0	0
			63	63		
7	F	50	Total	O	0	0
			50	50		
7	G	1	Total	O	0	0
			1	1		
7	H	2	Total	O	0	0
			2	2		

3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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.

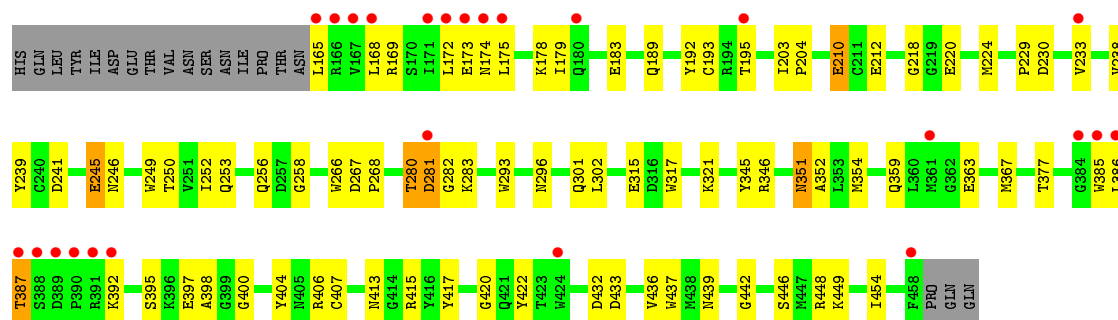
- Molecule 1: Fibrinogen alpha/alpha-E chain



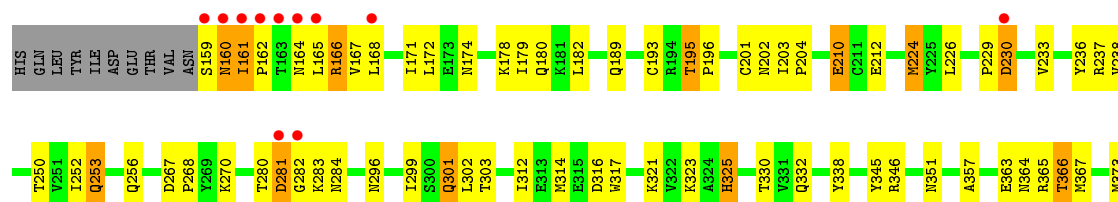
- Molecule 1: Fibrinogen alpha/alpha-E chain

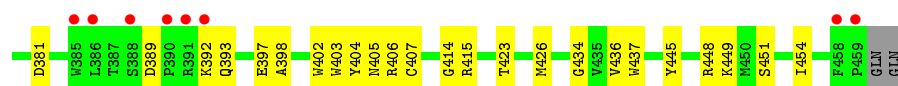


- Molecule 2: Fibrinogen beta chain

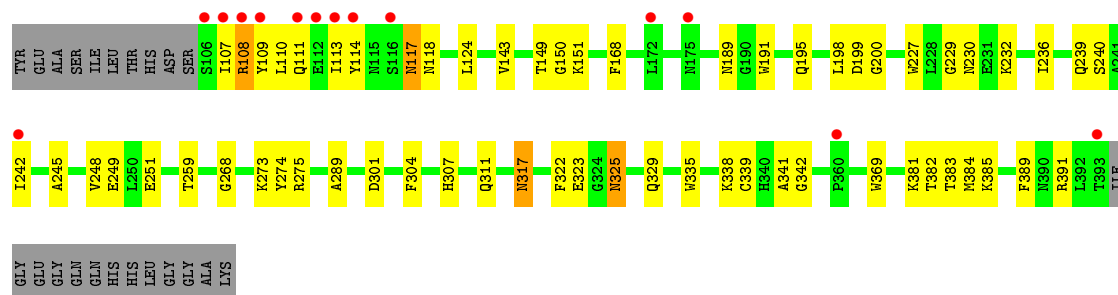
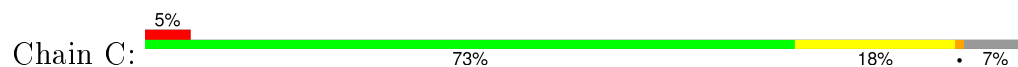


- Molecule 2: Fibrinogen beta chain

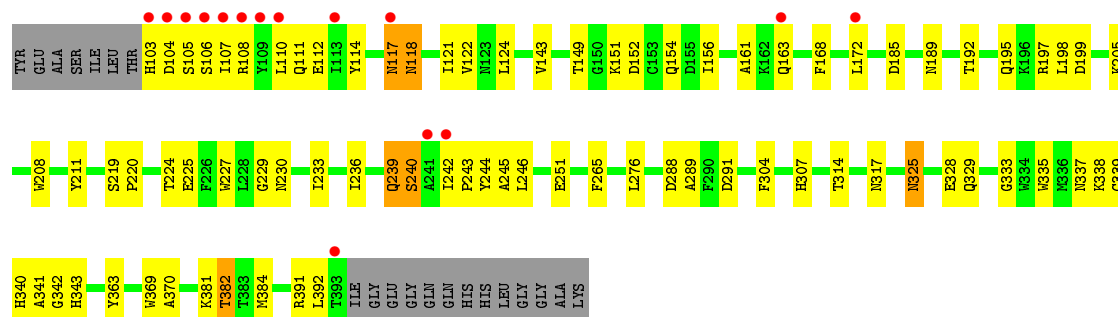




• Molecule 3: Fibrinogen gamma chain



• Molecule 3: Fibrinogen gamma chain



• Molecule 4: GHRP peptide



• Molecule 4: GHRP peptide



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	54.51Å 147.66Å 229.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	17.98 – 2.45 17.98 – 2.45	Depositor EDS
% Data completeness (in resolution range)	98.3 (17.98-2.45) 98.3 (17.98-2.45)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.10	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.53 (at 2.46Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.222 , 0.263 0.223 , 0.264	Depositor DCC
R_{free} test set	3443 reflections (5.06%)	DCC
Wilson B-factor (Å ²)	35.1	Xtriage
Anisotropy	0.086	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 38.2	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	1 of 68004 reflections (0.001%)	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	10794	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.98% 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, 0.375 respectively for untwinned datasets, and 0.333, 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: CA, NAG

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	0.28	0/491	0.52	0/655
1	D	0.33	0/517	0.56	0/687
2	B	0.38	0/2419	0.62	0/3264
2	E	0.37	0/2472	0.63	0/3339
3	C	0.39	0/2376	0.62	0/3213
3	F	0.38	0/2401	0.61	0/3247
4	G	0.45	0/34	0.60	0/43
4	H	0.48	0/34	0.53	0/43
All	All	0.37	0/10744	0.61	0/14491

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	490	0	508	40	0
1	D	516	0	543	41	0
2	B	2359	0	2232	74	0
2	E	2410	0	2281	88	0
3	C	2312	0	2165	64	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	F	2336	0	2181	75	0
4	G	33	0	32	7	0
4	H	33	0	32	9	0
5	B	28	0	25	1	0
5	E	28	0	25	1	0
6	B	1	0	0	0	0
6	C	2	0	0	0	0
6	E	1	0	0	0	0
6	F	1	0	0	0	0
7	A	2	0	0	0	0
7	B	63	0	0	2	0
7	C	54	0	0	2	0
7	D	9	0	0	0	0
7	E	63	0	0	3	0
7	F	50	0	0	1	0
7	G	1	0	0	1	0
7	H	2	0	0	1	0
All	All	10794	0	10024	347	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 17.

All (347) 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)
3:F:172:LEU:HD12	3:F:172:LEU:H	1.20	1.06
3:C:329:GLN:NE2	4:G:3:ARG:HH11	1.64	0.95
2:E:166:ARG:HD3	2:E:166:ARG:H	1.39	0.88
3:C:117:ASN:HD22	3:C:118:ASN:N	1.74	0.84
3:C:108:ARG:HD2	3:C:108:ARG:H	1.43	0.84
1:A:169:LEU:H	2:B:189:GLN:HE22	1.25	0.83
2:E:179:ILE:HD12	3:F:117:ASN:HB2	1.60	0.83
3:F:329:GLN:NE2	4:H:3:ARG:HH11	1.77	0.82
3:F:307:HIS:HE1	3:F:341:ALA:H	1.28	0.81
1:A:175:LEU:HD23	1:A:175:LEU:H	1.47	0.79
3:C:149:THR:HG23	3:C:168:PHE:O	1.83	0.79
3:F:149:THR:HG22	3:F:168:PHE:O	1.83	0.78
3:F:172:LEU:CD1	3:F:172:LEU:H	1.97	0.77
1:D:133:ILE:HD12	1:D:133:ILE:H	1.50	0.75
2:E:423:THR:H	2:E:426:MET:HE3	1.52	0.74
2:E:423:THR:N	2:E:426:MET:HE3	2.03	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:281:ASP:OD1	2:E:282:GLY:N	2.20	0.73
2:E:230:ASP:O	2:E:233:VAL:HG22	1.89	0.72
3:F:219:SER:OG	3:F:224:THR:HG22	1.90	0.72
3:F:329:GLN:HE21	4:H:3:ARG:NH1	1.89	0.71
1:D:129:LYS:O	1:D:129:LYS:HD2	1.91	0.71
2:E:252:ILE:HD13	2:E:454:ILE:HG12	1.73	0.71
1:A:153:ASP:O	1:A:157:LYS:HG2	1.91	0.71
2:E:168:LEU:O	2:E:172:LEU:HD23	1.91	0.70
3:C:108:ARG:HD2	3:C:108:ARG:N	2.07	0.69
1:D:176:LYS:HA	1:D:176:LYS:HE2	1.74	0.69
1:A:178:TYR:O	1:A:182:GLN:HG2	1.93	0.69
3:F:239:GLN:O	3:F:240:SER:HB3	1.92	0.69
3:C:329:GLN:HE21	4:G:3:ARG:HH11	1.41	0.69
2:E:161:ILE:H	2:E:161:ILE:HD13	1.58	0.68
3:C:151:LYS:HB3	3:C:239:GLN:HE22	1.57	0.68
3:C:240:SER:OG	3:C:242:ILE:HG12	1.94	0.68
3:F:195:GLN:OE1	3:F:382:THR:HG22	1.92	0.68
3:C:149:THR:HG21	7:C:433:HOH:O	1.94	0.68
2:B:230:ASP:O	2:B:233:VAL:HG22	1.93	0.68
2:B:210:GLU:OE1	2:B:212:GLU:HB3	1.93	0.68
3:F:172:LEU:N	3:F:172:LEU:HD12	2.02	0.67
3:F:329:GLN:NE2	4:H:3:ARG:NH1	2.40	0.67
3:F:307:HIS:CE1	3:F:341:ALA:H	2.12	0.67
1:D:139:ASN:HB3	3:F:114:TYR:CE2	2.29	0.67
3:C:149:THR:HG22	3:C:150:GLY:N	2.09	0.67
1:D:133:ILE:HD12	1:D:133:ILE:N	2.09	0.67
2:E:179:ILE:CD1	3:F:117:ASN:HB2	2.24	0.67
3:C:227:TRP:HZ2	3:C:230:ASN:HD21	1.40	0.67
2:B:386:LEU:O	2:B:387:THR:HB	1.94	0.66
3:C:195:GLN:OE1	3:C:382:THR:HG22	1.95	0.66
3:C:307:HIS:HE1	3:C:342:GLY:H	1.44	0.66
1:D:168:ALA:HA	2:E:189:GLN:HE22	1.60	0.65
3:C:329:GLN:NE2	4:G:3:ARG:NH1	2.41	0.65
2:B:385:TRP:HA	2:B:406:ARG:HD3	1.79	0.65
1:A:131:GLN:HE21	1:A:132:HIS:HB2	1.63	0.64
1:A:181:GLN:HE21	2:B:174:ASN:ND2	1.95	0.64
2:E:166:ARG:CD	2:E:166:ARG:H	2.11	0.63
2:E:267:ASP:HB3	2:E:268:PRO:HD3	1.80	0.63
2:E:167:VAL:O	2:E:171:ILE:HG12	1.99	0.63
2:B:229:PRO:HD2	2:B:233:VAL:HG21	1.80	0.63
1:A:181:GLN:HE21	2:B:174:ASN:HD22	1.46	0.63

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:161:ILE:N	2:E:162:PRO:HD2	2.13	0.62
2:E:373:MET:HE2	2:E:404:TYR:O	2.00	0.62
1:A:175:LEU:CD2	1:A:175:LEU:H	2.13	0.62
2:B:363:GLU:O	2:B:367:MET:HG2	1.99	0.62
1:D:137:GLN:NE2	1:D:189:ILE:HG23	2.14	0.62
1:D:144:LEU:HD13	1:D:182:GLN:HB3	1.82	0.62
2:E:448:ARG:HH11	2:E:448:ARG:HG2	1.65	0.62
1:A:166:SER:HB3	2:B:195:THR:HG23	1.82	0.61
1:A:139:ASN:HB3	3:C:114:TYR:CE2	2.35	0.61
3:F:329:GLN:HE21	4:H:3:ARG:HH11	1.43	0.61
3:C:329:GLN:HE22	4:G:3:ARG:HH11	1.44	0.61
1:D:169:LEU:H	2:E:189:GLN:NE2	1.98	0.61
3:C:227:TRP:HZ2	3:C:230:ASN:ND2	1.98	0.61
3:C:329:GLN:HE21	4:G:3:ARG:NH1	1.98	0.60
2:B:359:GLN:NE2	2:B:442:GLY:HA2	2.16	0.60
2:E:389:ASP:HB3	2:E:392:LYS:HD2	1.84	0.60
2:E:364:ASN:HA	2:E:367:MET:HE3	1.84	0.60
3:C:249:GLU:HB2	3:C:383:THR:HB	1.84	0.60
1:D:136:LEU:O	1:D:140:VAL:HG13	2.02	0.59
2:B:363:GLU:HG2	2:B:367:MET:CE	2.32	0.59
2:E:363:GLU:HA	2:E:366:THR:HG23	1.84	0.59
1:A:158:ILE:HG23	2:B:189:GLN:HE21	1.68	0.59
2:B:165:LEU:HD12	2:B:165:LEU:N	2.18	0.59
3:C:109:TYR:HD1	3:C:110:LEU:HD12	1.67	0.59
3:C:307:HIS:HE1	3:C:341:ALA:H	1.51	0.59
1:D:139:ASN:HB3	3:F:114:TYR:CZ	2.37	0.59
2:B:238:VAL:HG21	2:B:250:THR:HG23	1.84	0.58
3:C:323:GLU:CD	3:C:323:GLU:H	2.05	0.58
3:F:325:ASN:HD22	3:F:325:ASN:C	2.06	0.58
1:D:189:ILE:O	1:D:189:ILE:HG22	2.04	0.58
1:A:131:GLN:NE2	1:A:132:HIS:H	2.01	0.58
3:C:149:THR:HG22	3:C:150:GLY:H	1.68	0.57
3:F:185:ASP:OD2	3:F:189:ASN:HB2	2.03	0.57
2:B:174:ASN:O	2:B:178:LYS:HG2	2.04	0.57
3:C:338:LYS:N	3:C:339:CYS:HA	2.18	0.57
1:A:144:LEU:HD23	2:B:175:LEU:HD21	1.87	0.57
1:D:169:LEU:H	2:E:189:GLN:HE22	1.52	0.57
1:A:185:LEU:O	1:A:189:ILE:HG13	2.05	0.57
2:B:230:ASP:HB3	2:B:233:VAL:HG13	1.87	0.57
2:E:238:VAL:HG21	2:E:250:THR:HG23	1.86	0.57
3:F:117:ASN:O	3:F:121:ILE:HG13	2.05	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:246:LEU:HD22	3:F:265:PHE:CE1	2.40	0.56
2:B:245:GLU:O	2:B:246:ASN:HB2	2.05	0.56
3:C:117:ASN:C	3:C:117:ASN:HD22	2.06	0.56
1:A:133:ILE:HA	1:A:136:LEU:HB3	1.87	0.56
1:D:188:VAL:HG11	2:E:165:LEU:HG	1.87	0.56
5:B:470:NAG:H62	5:B:471:NAG:O5	2.06	0.56
2:E:159:SER:HA	2:E:162:PRO:HG3	1.87	0.56
1:A:140:VAL:HG21	2:B:168:LEU:HD11	1.87	0.56
2:E:280:THR:O	2:E:281:ASP:C	2.45	0.56
3:C:307:HIS:CE1	3:C:342:GLY:H	2.22	0.56
2:E:174:ASN:HD21	2:E:178:LYS:NZ	2.03	0.55
2:B:183:GLU:HG2	3:C:124:LEU:HD13	1.87	0.55
2:B:267:ASP:HB3	2:B:268:PRO:HD3	1.88	0.55
3:F:195:GLN:HE22	3:F:382:THR:HG21	1.70	0.55
2:B:302:LEU:HD13	2:B:454:ILE:HD11	1.87	0.55
2:B:406:ARG:N	2:B:407:CYS:HA	2.22	0.55
3:C:307:HIS:CE1	3:C:341:ALA:H	2.25	0.55
1:A:133:ILE:HD12	2:B:165:LEU:HD22	1.88	0.55
1:A:136:LEU:HD13	1:A:136:LEU:O	2.05	0.55
3:F:338:LYS:N	3:F:339:CYS:HA	2.21	0.55
3:C:242:ILE:O	3:C:242:ILE:HG13	2.07	0.55
2:E:389:ASP:OD2	2:E:392:LYS:HE3	2.07	0.55
3:F:251:GLU:HB3	3:F:381:LYS:HB2	1.88	0.55
3:F:117:ASN:HD22	3:F:118:ASN:N	2.05	0.54
1:A:140:VAL:HG11	2:B:172:LEU:HD21	1.89	0.54
3:C:189:ASN:OD1	3:C:391:ARG:HG3	2.07	0.54
3:F:199:ASP:O	3:F:225:GLU:OE2	2.26	0.54
1:D:133:ILE:H	1:D:133:ILE:CD1	2.17	0.54
3:F:104:ASP:C	3:F:106:SER:H	2.10	0.54
2:E:162:PRO:HA	2:E:165:LEU:HD13	1.88	0.54
2:B:315:GLU:HB3	2:B:449:LYS:HB2	1.88	0.54
3:C:200:GLY:HA2	7:C:419:HOH:O	2.07	0.54
3:F:151:LYS:HB3	3:F:239:GLN:HE22	1.73	0.54
3:F:325:ASN:ND2	3:F:328:GLU:H	2.06	0.54
2:E:224:MET:CE	2:E:237:ARG:HD3	2.38	0.54
1:D:185:LEU:O	1:D:189:ILE:HG13	2.09	0.53
1:D:135:LEU:O	1:D:135:LEU:HD13	2.09	0.53
1:D:165:CYS:HB3	2:E:193:CYS:HA	1.89	0.53
2:E:317:TRP:CE3	2:E:448:ARG:HD3	2.44	0.53
2:B:238:VAL:HG21	2:B:250:THR:CG2	2.39	0.53
1:D:136:LEU:HD21	3:F:111:GLN:HG2	1.91	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:117:ASN:HD22	3:F:117:ASN:C	2.12	0.53
3:F:195:GLN:OE1	3:F:382:THR:CG2	2.56	0.52
3:C:107:ILE:HD12	3:C:108:ARG:NH2	2.24	0.52
3:F:211:TYR:CE2	3:F:333:GLY:HA3	2.45	0.52
2:E:323:LYS:NZ	2:E:325:HIS:HD2	2.08	0.52
2:E:230:ASP:OD2	2:E:233:VAL:HG13	2.09	0.52
2:E:161:ILE:N	2:E:162:PRO:CD	2.71	0.52
3:F:151:LYS:HD2	3:F:172:LEU:HD11	1.91	0.51
2:E:373:MET:CE	2:E:405:ASN:HA	2.40	0.51
3:C:251:GLU:HB3	3:C:381:LYS:HB2	1.92	0.51
2:E:162:PRO:O	2:E:165:LEU:HB2	2.10	0.51
3:C:322:PHE:CZ	4:G:3:ARG:HG2	2.46	0.51
1:A:136:LEU:C	1:A:136:LEU:HD13	2.31	0.51
2:E:160:ASN:HD22	2:E:160:ASN:N	2.08	0.51
3:C:195:GLN:HB3	3:C:384:MET:HB2	1.92	0.51
2:B:241:ASP:HB3	2:B:249:TRP:HB2	1.93	0.51
2:B:169:ARG:O	2:B:173:GLU:HG2	2.11	0.51
2:B:172:LEU:HB3	3:C:113:ILE:HG22	1.93	0.51
2:E:201:CYS:O	3:F:143:VAL:HG21	2.11	0.51
2:E:229:PRO:CB	2:E:301:GLN:HE22	2.24	0.50
1:A:167:ARG:HD3	2:B:192:TYR:CE1	2.46	0.50
3:F:118:ASN:O	3:F:122:VAL:HG23	2.11	0.50
1:A:175:LEU:HD23	1:A:175:LEU:N	2.22	0.50
2:B:252:ILE:HD13	2:B:454:ILE:HG12	1.94	0.50
1:D:135:LEU:C	1:D:135:LEU:HD13	2.31	0.50
3:F:288:ASP:OD2	3:F:291:ASP:HB2	2.11	0.50
2:E:406:ARG:N	2:E:407:CYS:HA	2.27	0.49
2:E:161:ILE:CD1	2:E:161:ILE:H	2.23	0.49
1:D:158:ILE:HG23	2:E:189:GLN:HE21	1.78	0.49
2:B:363:GLU:HG2	2:B:367:MET:HE3	1.93	0.49
3:F:163:GLN:H	3:F:163:GLN:CD	2.15	0.49
2:E:270:LYS:HA	2:E:296:ASN:HB2	1.94	0.49
1:A:169:LEU:H	2:B:189:GLN:NE2	2.01	0.49
2:B:359:GLN:HE22	2:B:442:GLY:HA2	1.76	0.49
1:D:135:LEU:HD13	1:D:139:ASN:ND2	2.27	0.49
2:E:160:ASN:C	2:E:162:PRO:HD2	2.33	0.49
1:D:131:GLN:CD	1:D:131:GLN:N	2.66	0.49
3:F:245:ALA:HB1	3:F:392:LEU:HD11	1.94	0.49
4:G:3:ARG:NH2	7:G:40:HOH:O	2.41	0.49
1:D:133:ILE:O	1:D:137:GLN:HG2	2.13	0.49
3:F:172:LEU:CD1	3:F:239:GLN:HE21	2.26	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:205:LYS:HD2	7:F:439:HOH:O	2.13	0.48
2:B:398:ALA:O	2:B:415:ARG:HD3	2.13	0.48
2:E:398:ALA:O	2:E:415:ARG:HD3	2.13	0.48
2:E:332:GLN:O	2:E:338:TYR:HA	2.13	0.48
2:B:218:GLY:O	2:B:220:GLU:HG3	2.13	0.48
2:B:385:TRP:CD1	2:B:406:ARG:HD2	2.48	0.48
3:F:208:TRP:HA	3:F:314:THR:HG21	1.95	0.48
2:B:204:PRO:HG2	2:B:224:MET:O	2.12	0.48
2:B:386:LEU:O	2:B:387:THR:CB	2.59	0.48
2:E:314:MET:HA	2:E:449:LYS:O	2.12	0.48
2:B:422:TYR:OH	2:B:432:ASP:HB2	2.14	0.48
3:F:108:ARG:O	3:F:112:GLU:HG3	2.14	0.48
2:E:302:LEU:HD13	2:E:454:ILE:HD11	1.95	0.48
1:D:129:LYS:HG2	3:F:107:ILE:HD12	1.95	0.48
2:E:168:LEU:HD12	2:E:168:LEU:H	1.79	0.48
3:C:307:HIS:HD2	3:C:335:TRP:O	1.95	0.48
3:F:307:HIS:HD2	3:F:335:TRP:O	1.97	0.47
2:E:281:ASP:OD1	2:E:283:LYS:HD3	2.13	0.47
3:F:104:ASP:C	3:F:106:SER:N	2.67	0.47
1:A:131:GLN:O	1:A:134:GLN:HG2	2.14	0.47
2:E:203:ILE:HA	2:E:204:PRO:HD3	1.80	0.47
3:F:103:HIS:CE1	3:F:105:SER:HB3	2.50	0.47
1:A:140:VAL:CG1	2:B:172:LEU:HD21	2.45	0.47
2:B:238:VAL:HG22	2:B:239:TYR:N	2.30	0.47
2:E:202:ASN:HD22	2:E:284:ASN:HB2	1.79	0.47
2:B:363:GLU:HG2	2:B:367:MET:HE2	1.96	0.47
3:C:323:GLU:CD	3:C:323:GLU:N	2.68	0.47
3:C:117:ASN:HD22	3:C:118:ASN:H	1.59	0.47
3:C:149:THR:CG2	3:C:150:GLY:N	2.77	0.47
2:E:414:GLY:HA3	2:E:434:GLY:O	2.15	0.47
2:B:351:ASN:ND2	2:B:354:MET:H	2.11	0.47
2:B:179:ILE:HD12	3:C:117:ASN:HB2	1.97	0.47
1:A:155:ASP:HB2	1:A:173:VAL:HG21	1.96	0.47
2:B:281:ASP:OD1	2:B:282:GLY:N	2.38	0.47
1:A:165:CYS:HB3	2:B:193:CYS:HA	1.97	0.46
2:E:325:HIS:O	2:E:345:TYR:HA	2.15	0.46
4:H:3:ARG:NH2	7:H:5:HOH:O	2.48	0.46
1:D:137:GLN:HE21	1:D:189:ILE:HG23	1.78	0.46
2:E:166:ARG:HD3	2:E:166:ARG:N	2.16	0.46
2:E:436:VAL:CG1	2:E:437:TRP:N	2.79	0.46
3:C:275:ARG:HA	3:C:311:GLN:HA	1.97	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:F:276:LEU:HD23	3:F:276:LEU:C	2.35	0.46
3:F:152:ASP:OD2	3:F:244:TYR:OH	2.33	0.46
2:B:436:VAL:CG1	2:B:437:TRP:N	2.78	0.46
2:E:210:GLU:OE1	2:E:212:GLU:HB3	2.14	0.46
3:F:104:ASP:HB2	3:F:107:ILE:HD12	1.97	0.46
2:E:330:THR:HG22	7:E:477:HOH:O	2.16	0.46
3:F:118:ASN:HA	3:F:118:ASN:HD22	1.53	0.46
7:E:474:HOH:O	3:F:143:VAL:CG2	2.63	0.46
2:B:203:ILE:HA	2:B:204:PRO:HD3	1.90	0.46
3:C:325:ASN:HD22	3:C:325:ASN:C	2.19	0.46
3:C:107:ILE:HB	3:C:108:ARG:NE	2.31	0.45
2:E:161:ILE:N	2:E:161:ILE:HD13	2.29	0.45
2:B:245:GLU:O	2:B:246:ASN:CB	2.65	0.45
3:C:289:ALA:HB3	3:C:369:TRP:CE2	2.52	0.45
2:E:161:ILE:HD13	2:E:162:PRO:HD3	1.99	0.45
2:B:175:LEU:O	2:B:178:LYS:HB2	2.17	0.45
2:B:280:THR:O	2:B:282:GLY:N	2.49	0.45
3:F:340:HIS:CE1	4:H:1:GLY:N	2.85	0.45
3:F:289:ALA:HB3	3:F:369:TRP:CE2	2.51	0.45
4:H:2:HIS:CD2	4:H:4:PRO:HD3	2.51	0.45
1:A:180:ASP:O	1:A:184:GLN:CB	2.64	0.45
2:E:357:ALA:O	2:E:365:ARG:HG3	2.17	0.45
1:A:188:VAL:HG12	1:A:188:VAL:O	2.17	0.45
2:E:381:ASP:HB2	2:E:393:GLN:HE21	1.82	0.45
1:D:166:SER:HB3	2:E:195:THR:HG23	1.99	0.45
3:C:117:ASN:C	3:C:117:ASN:ND2	2.69	0.44
2:E:160:ASN:N	2:E:160:ASN:ND2	2.64	0.44
3:F:340:HIS:ND1	3:F:343:HIS:HB2	2.32	0.44
3:C:198:LEU:HD12	3:C:199:ASP:N	2.32	0.44
3:F:195:GLN:HB3	3:F:384:MET:HB2	1.99	0.44
2:E:202:ASN:ND2	2:E:284:ASN:HB2	2.33	0.44
1:A:186:GLU:CD	1:A:189:ILE:HD12	2.38	0.44
1:D:131:GLN:CD	1:D:131:GLN:H	2.20	0.44
1:A:176:LYS:HA	1:A:179:GLU:OE1	2.17	0.44
3:F:363:TYR:HB3	4:H:3:ARG:NH2	2.32	0.44
3:F:197:ARG:HB2	3:F:382:THR:HB	1.99	0.44
1:D:179:GLU:O	1:D:183:LYS:HG3	2.17	0.44
2:B:258:GLY:HA2	7:B:500:HOH:O	2.17	0.44
2:E:159:SER:HA	2:E:162:PRO:CG	2.48	0.44
2:B:283:LYS:HD2	2:B:283:LYS:N	2.33	0.44
1:A:137:GLN:O	1:A:140:VAL:HG22	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:C:273:LYS:HE2	3:C:317:ASN:HD21	1.83	0.44
1:A:150:LEU:HD21	3:C:124:LEU:HD23	1.99	0.44
1:A:180:ASP:O	1:A:184:GLN:HB2	2.17	0.44
2:B:420:GLY:HA2	2:B:446:SER:O	2.18	0.44
2:E:299:ILE:O	2:E:303:THR:HG23	2.18	0.43
2:B:345:TYR:CG	2:B:346:ARG:N	2.86	0.43
3:F:307:HIS:CE1	3:F:342:GLY:H	2.37	0.43
2:B:397:GLU:O	2:B:398:ALA:HB3	2.18	0.43
1:D:132:HIS:O	1:D:136:LEU:HB2	2.18	0.43
3:C:117:ASN:ND2	3:C:118:ASN:N	2.56	0.43
2:B:385:TRP:CH2	2:B:392:LYS:HB3	2.53	0.43
1:A:139:ASN:HB3	3:C:114:TYR:CZ	2.53	0.43
1:D:178:TYR:O	1:D:182:GLN:HG2	2.19	0.43
1:A:186:GLU:C	1:A:188:VAL:H	2.23	0.43
2:B:293:TRP:HE1	2:B:296:ASN:ND2	2.17	0.43
3:F:227:TRP:HZ2	3:F:230:ASN:HD21	1.66	0.43
1:D:136:LEU:HD22	1:D:136:LEU:HA	1.84	0.43
3:C:236:ILE:O	3:C:239:GLN:HG2	2.18	0.43
7:E:530:HOH:O	3:F:220:PRO:HD2	2.18	0.43
1:D:140:VAL:HG11	2:E:172:LEU:HD21	1.99	0.42
2:E:373:MET:HE1	2:E:405:ASN:HA	2.00	0.42
3:C:389:PHE:C	3:C:391:ARG:H	2.22	0.42
2:B:351:ASN:C	2:B:351:ASN:HD22	2.22	0.42
2:E:253:GLN:NE2	2:E:451:SER:HA	2.34	0.42
1:D:151:GLU:HG2	1:D:173:VAL:HG13	2.01	0.42
1:D:181:GLN:OE1	2:E:174:ASN:ND2	2.53	0.42
3:C:227:TRP:CZ2	3:C:230:ASN:ND2	2.85	0.42
1:D:151:GLU:OE2	2:E:182:LEU:HD21	2.20	0.42
2:B:395:SER:HB2	2:B:404:TYR:CE2	2.54	0.42
2:B:367:MET:HB2	2:B:406:ARG:HB3	2.01	0.42
3:C:113:ILE:HG22	3:C:113:ILE:O	2.18	0.42
3:F:192:THR:HG21	3:F:236:ILE:HD13	2.00	0.42
3:C:108:ARG:CD	3:C:108:ARG:N	2.73	0.42
1:D:185:LEU:HG	1:D:189:ILE:HD11	2.01	0.42
2:E:397:GLU:O	2:E:398:ALA:HB3	2.19	0.42
1:A:151:GLU:CG	1:A:173:VAL:HG13	2.50	0.42
2:E:316:ASP:HB2	2:E:445:TYR:OH	2.19	0.42
2:E:226:LEU:HD12	2:E:236:TYR:C	2.40	0.42
3:F:340:HIS:CE1	3:F:343:HIS:HB2	2.54	0.42
3:F:189:ASN:OD1	3:F:391:ARG:HG3	2.20	0.42
2:B:280:THR:O	2:B:281:ASP:C	2.57	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:174:ASN:HD21	2:E:178:LYS:HZ3	1.66	0.42
3:C:268:GLY:O	3:C:274:TYR:HA	2.20	0.42
5:E:471:NAG:H3	5:E:471:NAG:H83	2.02	0.42
2:E:229:PRO:HB3	2:E:301:GLN:HE22	1.84	0.41
3:F:105:SER:HA	3:F:108:ARG:NE	2.35	0.41
3:F:172:LEU:HD13	3:F:239:GLN:HB2	2.01	0.41
3:F:329:GLN:HE22	4:H:3:ARG:HH11	1.63	0.41
3:C:191:TRP:CE3	3:C:385:LYS:HG3	2.55	0.41
2:E:312:ILE:HA	2:E:451:SER:O	2.21	0.41
2:E:389:ASP:HB3	2:E:392:LYS:CD	2.50	0.41
2:B:317:TRP:CE3	2:B:448:ARG:HD3	2.56	0.41
3:C:229:GLY:H	3:C:232:LYS:HD2	1.85	0.41
3:C:248:VAL:O	3:C:259:THR:HA	2.20	0.41
1:A:144:LEU:HD23	2:B:175:LEU:HD11	2.03	0.41
2:E:367:MET:HB2	2:E:406:ARG:HB2	2.03	0.41
2:E:345:TYR:CG	2:E:346:ARG:N	2.88	0.41
2:E:402:TRP:CG	2:E:403:TRP:N	2.89	0.41
2:B:266:TRP:HA	2:B:377:THR:HG21	2.02	0.41
3:F:106:SER:O	3:F:110:LEU:HD13	2.21	0.41
2:E:167:VAL:HG22	2:E:168:LEU:HD12	2.03	0.41
3:C:230:ASN:HD22	3:C:230:ASN:HA	1.67	0.41
2:E:174:ASN:ND2	2:E:178:LYS:HZ3	2.19	0.41
3:F:143:VAL:HG23	3:F:143:VAL:O	2.21	0.41
1:D:151:GLU:CG	1:D:173:VAL:HG13	2.51	0.41
3:F:304:PHE:O	3:F:337:ASN:HB3	2.20	0.41
3:F:242:ILE:HA	3:F:243:PRO:HD2	1.95	0.41
2:B:400:GLY:HA3	2:B:413:ASN:O	2.21	0.41
1:D:140:VAL:CG1	2:E:172:LEU:HD21	2.51	0.41
7:B:473:HOH:O	3:C:143:VAL:HG22	2.21	0.41
2:B:417:TYR:OH	2:B:433:ASP:OD1	2.36	0.41
2:E:448:ARG:NH1	2:E:448:ARG:HG2	2.33	0.40
3:C:245:ALA:HB2	3:C:389:PHE:HD1	1.86	0.40
3:F:152:ASP:OD1	3:F:154:GLN:N	2.55	0.40
1:D:150:LEU:HD21	3:F:124:LEU:HD23	2.02	0.40
2:E:238:VAL:HG21	2:E:250:THR:CG2	2.50	0.40
2:B:229:PRO:HD2	2:B:233:VAL:CG2	2.49	0.40
3:C:304:PHE:CD1	3:C:338:LYS:HE3	2.57	0.40
1:A:186:GLU:OE1	1:A:189:ILE:HD12	2.22	0.40
1:D:134:GLN:H	1:D:134:GLN:HG2	1.74	0.40
3:F:229:GLY:O	3:F:233:ILE:HG13	2.21	0.40
2:B:230:ASP:H	2:B:233:VAL:HG22	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:166:SER:CB	2:B:195:THR:HG23	2.49	0.40
3:F:156:ILE:HG22	3:F:161:ALA:CB	2.51	0.40
2:B:352:ALA:HB2	2:B:439:ASN:ND2	2.37	0.40

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	58/66 (88%)	53 (91%)	5 (9%)	0	100	100
1	D	61/66 (92%)	58 (95%)	2 (3%)	1 (2%)	12	11
2	B	292/313 (93%)	272 (93%)	17 (6%)	3 (1%)	19	21
2	E	299/313 (96%)	278 (93%)	18 (6%)	3 (1%)	19	21
3	C	286/311 (92%)	271 (95%)	15 (5%)	0	100	100
3	F	289/311 (93%)	272 (94%)	15 (5%)	2 (1%)	26	32
4	G	2/4 (50%)	2 (100%)	0	0	100	100
4	H	2/4 (50%)	2 (100%)	0	0	100	100
All	All	1289/1388 (93%)	1208 (94%)	72 (6%)	9 (1%)	26	32

All (9) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	281	ASP
2	B	387	THR
2	E	281	ASP
3	F	240	SER
2	B	256	GLN
2	E	160	ASN
2	E	256	GLN

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Mol	Chain	Res	Type
3	F	370	ALA
1	D	189	ILE

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	55/61 (90%)	53 (96%)	2 (4%)	42	58
1	D	58/61 (95%)	55 (95%)	3 (5%)	29	40
2	B	251/270 (93%)	244 (97%)	7 (3%)	51	68
2	E	258/270 (96%)	243 (94%)	15 (6%)	25	34
3	C	242/259 (93%)	236 (98%)	6 (2%)	55	72
3	F	245/259 (95%)	238 (97%)	7 (3%)	50	66
4	G	3/3 (100%)	2 (67%)	1 (33%)	0	0
4	H	3/3 (100%)	3 (100%)	0	100	100
All	All	1115/1186 (94%)	1074 (96%)	41 (4%)	41	57

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

Mol	Chain	Res	Type
1	A	135	LEU
1	A	175	LEU
2	B	210	GLU
2	B	245	GLU
2	B	253	GLN
2	B	280	THR
2	B	301	GLN
2	B	321	LYS
2	B	351	ASN
3	C	108	ARG
3	C	111	GLN
3	C	117	ASN
3	C	301	ASP

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Mol	Chain	Res	Type
3	C	317	ASN
3	C	325	ASN
1	D	134	GLN
1	D	136	LEU
1	D	176	LYS
2	E	161	ILE
2	E	164	ASN
2	E	166	ARG
2	E	180	GLN
2	E	195	THR
2	E	196	PRO
2	E	210	GLU
2	E	224	MET
2	E	230	ASP
2	E	253	GLN
2	E	301	GLN
2	E	321	LYS
2	E	325	HIS
2	E	351	ASN
2	E	366	THR
3	F	117	ASN
3	F	118	ASN
3	F	198	LEU
3	F	239	GLN
3	F	317	ASN
3	F	325	ASN
3	F	382	THR
4	G	3	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (66) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	131	GLN
1	A	134	GLN
1	A	139	ASN
1	A	143	GLN
1	A	181	GLN
1	A	182	GLN
2	B	180	GLN
2	B	189	GLN
2	B	253	GLN
2	B	256	GLN

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Mol	Chain	Res	Type
2	B	296	ASN
2	B	301	GLN
2	B	332	GLN
2	B	339	GLN
2	B	351	ASN
2	B	439	ASN
3	C	117	ASN
3	C	118	ASN
3	C	130	GLN
3	C	230	ASN
3	C	239	GLN
3	C	254	ASN
3	C	307	HIS
3	C	317	ASN
3	C	319	ASN
3	C	325	ASN
3	C	329	GLN
3	C	350	GLN
1	D	137	GLN
1	D	139	ASN
1	D	143	GLN
1	D	181	GLN
1	D	182	GLN
1	D	184	GLN
2	E	160	ASN
2	E	164	ASN
2	E	174	ASN
2	E	180	GLN
2	E	189	GLN
2	E	202	ASN
2	E	228	GLN
2	E	253	GLN
2	E	256	GLN
2	E	271	GLN
2	E	296	ASN
2	E	301	GLN
2	E	325	HIS
2	E	332	GLN
2	E	339	GLN
2	E	351	ASN
2	E	393	GLN
2	E	439	ASN

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Mol	Chain	Res	Type
3	F	103	HIS
3	F	117	ASN
3	F	118	ASN
3	F	130	GLN
3	F	177	GLN
3	F	230	ASN
3	F	239	GLN
3	F	254	ASN
3	F	307	HIS
3	F	317	ASN
3	F	319	ASN
3	F	325	ASN
3	F	329	GLN
4	G	2	HIS

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 ⓘ

4 carbohydrates 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	$\# Z > 2$
5	NAG	B	470	2,5	14,14,15	0.71	0	15,19,21	1.23	2 (13%)
5	NAG	B	471	5	14,14,15	0.58	0	15,19,21	0.72	1 (6%)
5	NAG	E	470	2,5	14,14,15	0.61	0	15,19,21	0.76	1 (6%)
5	NAG	E	471	5	14,14,15	0.64	0	15,19,21	0.59	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
5	NAG	B	470	2,5	-	0/6/23/26	0/1/1/1
5	NAG	B	471	5	-	0/6/23/26	0/1/1/1
5	NAG	E	470	2,5	-	0/6/23/26	0/1/1/1
5	NAG	E	471	5	-	0/6/23/26	0/1/1/1

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	B	471	NAG	C2-N2-C7	-2.26	120.13	123.04
5	E	470	NAG	C2-N2-C7	-2.09	120.35	123.04
5	B	470	NAG	C3-C4-C5	2.58	114.70	110.20
5	B	470	NAG	C4-C3-C2	2.61	115.28	111.23

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	470	NAG	1	0
5	B	471	NAG	1	0
5	E	471	NAG	1	0

5.6 Ligand geometry

Of 5 ligands modelled in this entry, 5 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers

There are no such residues in this entry.

5.8 Polymer linkage issues

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	60/66 (90%)	1.29	15 (25%) 1 1	32, 73, 110, 113	0
1	D	63/66 (95%)	0.34	6 (9%) 10 10	27, 45, 93, 99	0
2	B	294/313 (93%)	0.09	25 (8%) 13 13	20, 33, 84, 114	0
2	E	301/313 (96%)	-0.00	19 (6%) 23 25	22, 33, 73, 112	0
3	C	288/311 (92%)	-0.09	14 (4%) 33 36	19, 30, 65, 115	0
3	F	291/311 (93%)	0.03	15 (5%) 31 34	21, 35, 58, 108	0
4	G	4/4 (100%)	0.61	0 100 100	33, 35, 46, 50	0
4	H	4/4 (100%)	0.45	0 100 100	25, 31, 34, 46	0
All	All	1305/1388 (94%)	0.09	94 (7%) 18 19	19, 34, 87, 115	0

All (94) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	166	ARG	8.4
2	B	391	ARG	6.9
1	D	129	LYS	6.5
2	E	159	SER	5.9
3	F	393	THR	5.8
2	E	459	PRO	5.4
1	A	190	ALA	5.4
1	D	130	VAL	5.2
2	E	160	ASN	4.7
2	E	163	THR	4.7
3	F	103	HIS	4.6
3	C	109	TYR	4.5
1	A	183	LYS	4.5
2	E	391	ARG	4.5
2	B	386	LEU	4.5
1	A	172	GLU	4.4

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Mol	Chain	Res	Type	RSRZ
2	B	167	VAL	4.3
3	C	108	ARG	4.3
2	B	392	LYS	4.3
1	A	135	LEU	4.2
1	A	189	ILE	4.2
3	C	112	GLU	4.1
2	B	385	TRP	4.1
2	E	386	LEU	3.8
2	E	161	ILE	3.8
1	A	134	GLN	3.7
2	E	281	ASP	3.7
3	F	108	ARG	3.5
2	B	390	PRO	3.4
3	C	107	ILE	3.4
2	B	173	GLU	3.3
2	B	195	THR	3.3
1	A	132	HIS	3.3
3	F	104	ASP	3.3
2	E	458	PHE	3.2
3	C	106	SER	3.2
2	E	164	ASN	3.2
2	B	174	ASN	3.1
2	B	458	PHE	3.1
1	A	131	GLN	3.0
3	F	242	ILE	3.0
1	A	187	GLN	3.0
1	A	133	ILE	3.0
2	B	168	LEU	2.9
2	E	388	SER	2.9
1	A	170	ALA	2.9
2	B	165	LEU	2.9
3	F	106	SER	2.8
2	E	230	ASP	2.8
3	F	109	TYR	2.7
2	B	281	ASP	2.7
2	E	165	LEU	2.7
3	C	172	LEU	2.7
3	F	110	LEU	2.7
2	B	233	VAL	2.7
2	B	361	MET	2.7
1	D	190	ALA	2.6
2	E	385	TRP	2.6

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Mol	Chain	Res	Type	RSRZ
2	E	390	PRO	2.6
2	B	180	GLN	2.6
2	B	172	LEU	2.5
2	B	175	LEU	2.5
2	E	162	PRO	2.5
3	C	175	ASN	2.5
3	F	113	ILE	2.5
3	C	113	ILE	2.5
2	B	424	TRP	2.4
1	A	136	LEU	2.4
3	F	117	ASN	2.4
2	B	389	ASP	2.4
1	A	182	GLN	2.4
2	E	282	GLY	2.4
3	C	393	THR	2.4
1	D	189	ILE	2.4
3	F	241	ALA	2.4
3	C	360	PRO	2.4
2	E	392	LYS	2.4
2	B	387	THR	2.3
2	E	168	LEU	2.3
1	A	138	LYS	2.3
2	B	388	SER	2.3
3	F	107	ILE	2.2
1	D	187	GLN	2.2
1	D	191	LYS	2.2
3	C	242	ILE	2.2
2	B	171	ILE	2.2
1	A	176	LYS	2.2
3	C	116	SER	2.1
3	F	105	SER	2.1
3	F	172	LEU	2.1
3	F	163	GLN	2.0
3	C	111	GLN	2.0
2	B	384	GLY	2.0
3	C	114	TYR	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](#)

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
5	NAG	E	470	14/15	0.81	0.32	-	56,61,66,74	0
5	NAG	B	471	14/15	0.51	0.63	-	110,112,113,113	0
5	NAG	E	471	14/15	0.72	0.59	-	79,84,87,88	0
5	NAG	B	470	14/15	0.74	0.44	-	93,96,99,105	0

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(Å ²)	Q<0.9
6	CA	C	408	1/1	0.99	0.10	-1.13	31,31,31,31	0
6	CA	C	407	1/1	0.94	0.05	-1.74	32,32,32,32	0
6	CA	F	407	1/1	0.99	0.03	-1.96	27,27,27,27	0
6	CA	E	472	1/1	0.89	0.05	-	46,46,46,46	0
6	CA	B	472	1/1	0.94	0.07	-	44,44,44,44	0

6.5 Other polymers [i](#)

There are no such residues in this entry.