



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 10:32 PM GMT

PDB ID : 4Y4F
Title : Crystal structure of the mCD1d/GCK127/iNKTCR ternary complex
Authors : Zajonc, D.M.; Yu, E.D.
Deposited on : 2015-02-10
Resolution : 3.19 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

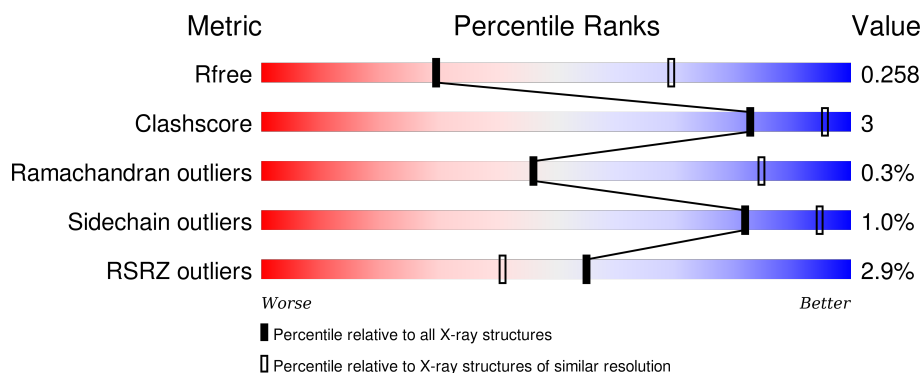
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.19 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.



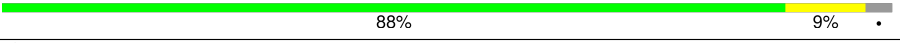


Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	91344	1124 (3.24-3.16)
Clashscore	102246	1024 (3.22-3.18)
Ramachandran outliers	100387	1004 (3.22-3.18)
Sidechain outliers	100360	1003 (3.22-3.18)
RSRZ outliers	91569	1129 (3.24-3.16)

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	285	<div> <div>5%</div> <div>85%</div> <div>9%</div> <div>6%</div> </div>
1	E	285	<div> <div>8%</div> <div>84%</div> <div>10%</div> <div>6%</div> </div>
2	B	99	<div> <div>4%</div> <div>92%</div> <div>5%</div> <div>•</div> </div>
2	F	99	<div> <div>•</div> <div>89%</div> <div>8%</div> <div>•</div> </div>
3	C	209	<div> <div>89%</div> <div>8%</div> <div>•</div> </div>

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Mol	Chain	Length	Quality of chain
3	G	209	
4	D	241	
4	H	241	

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
6	49M	A	504	-	-	-	X
6	49M	E	504	-	-	-	X

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 12597 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 Antigen-presenting glycoprotein CD1d1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	267	Total	C	N	O	S	0	0	0
			2076	1331	345	387	13			
1	E	267	Total	C	N	O	S	0	0	0
			2025	1292	337	383	13			

There are 14 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	201	HIS	ASP	variant	UNP P11609
A	280	HIS	-	expression tag	UNP P11609
A	281	HIS	-	expression tag	UNP P11609
A	282	HIS	-	expression tag	UNP P11609
A	283	HIS	-	expression tag	UNP P11609
A	284	HIS	-	expression tag	UNP P11609
A	285	HIS	-	expression tag	UNP P11609
E	201	HIS	ASP	variant	UNP P11609
E	280	HIS	-	expression tag	UNP P11609
E	281	HIS	-	expression tag	UNP P11609
E	282	HIS	-	expression tag	UNP P11609
E	283	HIS	-	expression tag	UNP P11609
E	284	HIS	-	expression tag	UNP P11609
E	285	HIS	-	expression tag	UNP P11609

- Molecule 2 is a protein called Beta-2-microglobulin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	B	96	Total	C	N	O	S	0	0	0
			755	483	125	141	6			
2	F	96	Total	C	N	O	S	0	0	0
			753	481	127	139	6			

- Molecule 3 is a protein called Chimeric TCR Valpha14/Jalpha18 chain (mouse variable do-

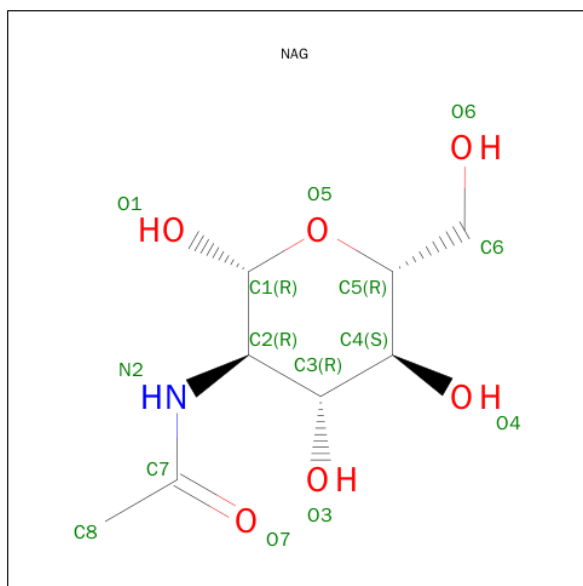
main/ human constant domain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
3	C	202	Total	C	N	O	S	0	1	0
			1537	951	262	316	8			
3	G	202	Total	C	N	O	S	0	0	0
			1521	941	259	313	8			

- Molecule 4 is a protein called Chimeric TCR Vbeta8.2 chain (mouse variable domain, human constant domain).

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
4	D	239	Total	C	N	O	S	0	0	0
			1858	1165	331	356	6			
4	H	239	Total	C	N	O	S	0	0	0
			1862	1168	332	356	6			

- Molecule 5 is N-ACETYL-D-GLUCOSAMINE (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



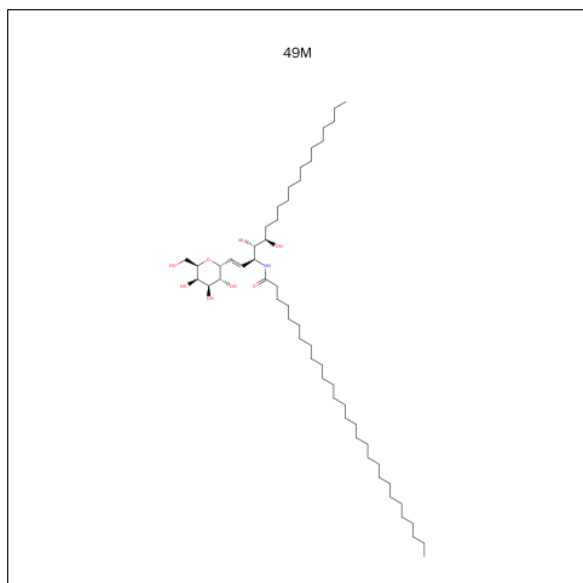
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	A	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
5	E	1	Total	C	N	O	0	0
			14	8	1	5		
5	E	1	Total	C	N	O	0	0
			14	8	1	5		

- Molecule 6 is (1R)-1,5-anhydro-1-[(1E,3S,4S,5R)-4,5-dihydroxy-3-(nonacosanoylamino)nona dec-1-en-1-yl]-D-galactitol (three-letter code: 49M) (formula: C₅₄H₁₀₅NO₈).

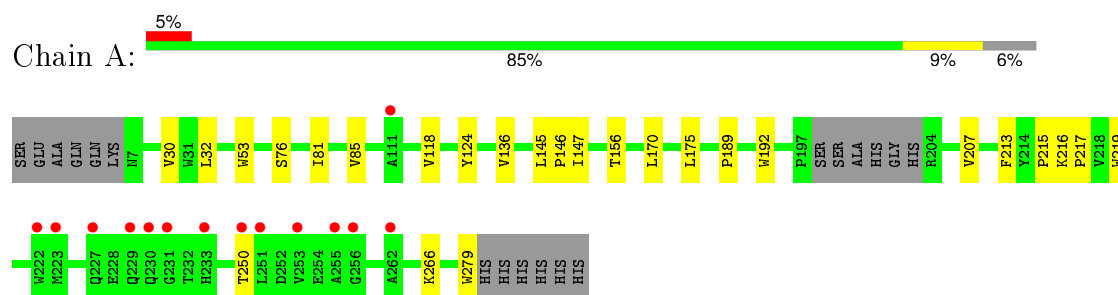


Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
6	A	1	Total	C	N	O	0	0
			63	54	1	8		
6	E	1	Total	C	N	O	0	0
			63	54	1	8		

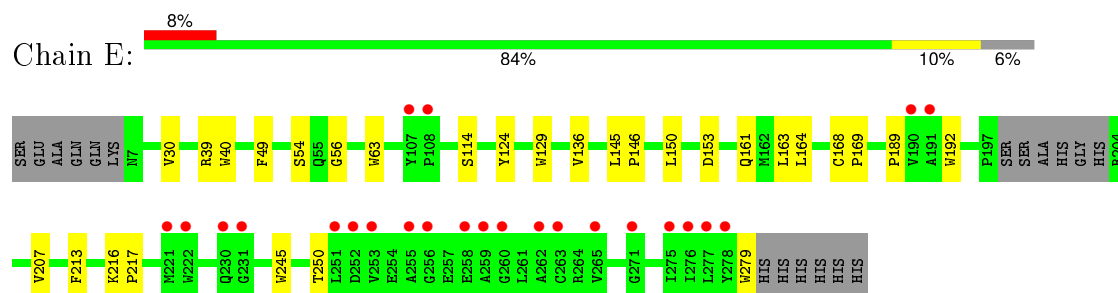
3 Residue-property plots [i](#)

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.

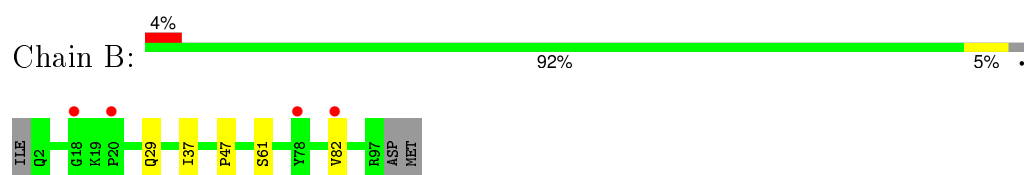
- Molecule 1: Antigen-presenting glycoprotein CD1d1



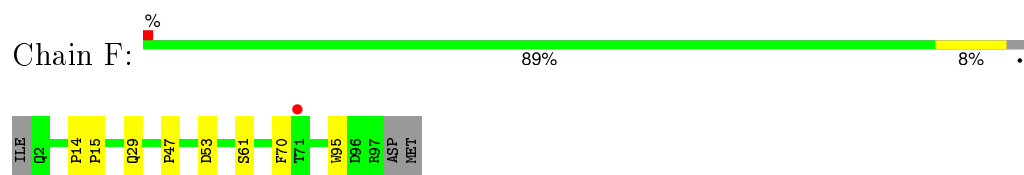
- Molecule 1: Antigen-presenting glycoprotein CD1d1




- Molecule 2: Beta-2-microglobulin



- Molecule 2: Beta-2-microglobulin




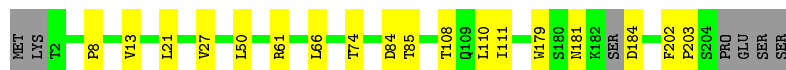
- Molecule 3: Chimeric TCR Valpha14/Jalpha18 chain (mouse variable domain/ human constant domain)

Chain C:  89% 8% .




- Molecule 3: Chimeric TCR Valpha14/Jalpha18 chain (mouse variable domain/ human constant domain)

Chain G:  88% 9% .




- Molecule 4: Chimeric TCR Vbeta8.2 chain (mouse variable domain, human constant domain)

Chain D:  88% 11% .



- Molecule 4: Chimeric TCR Vbeta8.2 chain (mouse variable domain, human constant domain)

Chain H:  89% 10% ..



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	78.59Å 149.73Å 101.39Å 90.00° 96.46° 90.00°	Depositor
Resolution (Å)	35.09 – 3.19 35.09 – 3.19	Depositor EDS
% Data completeness (in resolution range)	98.5 (35.09-3.19) 98.6 (35.09-3.19)	Depositor EDS
R_{merge}	0.13	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.90 (at 3.18Å)	Xtriage
Refinement program	REFMAC	Depositor
R, R_{free}	0.218 , 0.257 0.217 , 0.258	Depositor DCC
R_{free} test set	1190 reflections (3.21%)	DCC
Wilson B-factor (Å ²)	78.6	Xtriage
Anisotropy	0.027	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.29 , 34.3	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.31$	Xtriage
Outliers	1 of 38263 reflections (0.003%)	Xtriage
F_o, F_c correlation	0.92	EDS
Total number of atoms	12597	wwPDB-VP
Average B, all atoms (Å ²)	90.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.36% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 49M, 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	0.61	3/2138 (0.1%)	0.50	0/2919
1	E	0.62	6/2085 (0.3%)	0.49	0/2854
2	B	0.53	0/781	0.48	0/1070
2	F	0.53	1/779 (0.1%)	0.50	0/1068
3	C	0.43	1/1567 (0.1%)	0.48	0/2134
3	G	0.43	1/1548 (0.1%)	0.50	0/2111
4	D	0.51	3/1909 (0.2%)	0.49	0/2606
4	H	0.51	1/1913 (0.1%)	0.49	0/2610
All	All	0.53	16/12720 (0.1%)	0.49	0/17372

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	35	TRP	CD2-CE2	5.20	1.47	1.41
1	E	40	TRP	CD2-CE2	5.20	1.47	1.41
1	A	192	TRP	CD2-CE2	5.17	1.47	1.41
3	G	179	TRP	CD2-CE2	5.15	1.47	1.41
4	D	237	TRP	CD2-CE2	5.10	1.47	1.41
1	E	245	TRP	CD2-CE2	5.07	1.47	1.41
1	A	279	TRP	CD2-CE2	5.07	1.47	1.41
1	E	129	TRP	CD2-CE2	5.06	1.47	1.41
1	E	63	TRP	CD2-CE2	5.05	1.47	1.41
4	D	220	TRP	CD2-CE2	5.05	1.47	1.41
4	H	237	TRP	CD2-CE2	5.04	1.47	1.41
4	D	34	TRP	CD2-CE2	5.04	1.47	1.41
1	E	192	TRP	CD2-CE2	5.02	1.47	1.41
2	F	95	TRP	CD2-CE2	5.02	1.47	1.41
1	A	53	TRP	CD2-CE2	5.01	1.47	1.41
1	E	279	TRP	CD2-CE2	5.01	1.47	1.41

There are no bond angle outliers.

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	2076	0	1924	14	0
1	E	2025	0	1810	17	0
2	B	755	0	684	2	0
2	F	753	0	680	3	0
3	C	1537	0	1440	7	0
3	G	1521	0	1410	9	0
4	D	1858	0	1741	12	0
4	H	1862	0	1752	12	0
5	A	42	0	38	0	0
5	E	42	0	38	1	0
6	A	63	0	105	4	0
6	E	63	0	105	4	0
All	All	12597	0	11727	75	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 3.

All (75) 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)
4:H:115:ARG:HB2	4:H:222:GLN:HE22	1.50	0.76
4:D:204:HIS:NE2	4:D:235:GLU:HG3	2.01	0.74
1:E:49:PHE:CD1	1:E:54:SER:HB2	2.23	0.72
1:E:30:VAL:HG21	6:E:504:49M:H80	1.72	0.70
1:E:49:PHE:HD1	1:E:54:SER:HB2	1.59	0.67
1:A:219:TRP:HB3	1:A:266:LYS:HB2	1.78	0.66
1:E:189:PRO:HB3	1:E:213:PHE:HB3	1.79	0.64
1:A:30:VAL:HG21	6:A:504:49M:H80	1.82	0.61
2:F:29:GLN:HA	2:F:61:SER:HB2	1.84	0.60
1:E:145:LEU:HB3	1:E:146:PRO:HD3	1.82	0.60
3:C:85:THR:HG23	3:C:111:ILE:HA	1.84	0.59

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:85:THR:HG23	3:G:111:ILE:HA	1.85	0.57
1:A:216:LYS:N	1:A:217:PRO:HD2	2.21	0.56
1:E:207:VAL:HG22	1:E:250:THR:HG22	1.88	0.56
1:A:124:TYR:CZ	1:A:136:VAL:HG11	2.41	0.55
1:A:189:PRO:HB3	1:A:213:PHE:HB3	1.88	0.54
1:A:215:PRO:HB2	1:A:217:PRO:HD2	1.89	0.54
1:E:150:LEU:HD11	6:E:504:49M:H13	1.90	0.54
4:H:127:PRO:HD3	4:H:140:LEU:HG	1.89	0.54
4:H:150:ASP:HB2	4:H:173:PRO:HG2	1.91	0.53
1:A:81:ILE:O	1:A:85:VAL:HG23	2.10	0.52
4:D:199:GLN:HA	4:D:239:ARG:O	2.10	0.52
1:E:161:GLN:HB3	5:E:502:NAG:H82	1.93	0.51
2:B:29:GLN:HA	2:B:61:SER:HB2	1.91	0.51
4:D:124:LEU:HD11	4:D:140:LEU:HD23	1.92	0.51
1:A:156:THR:HG23	6:A:504:49M:H52	1.92	0.50
1:E:49:PHE:HD1	1:E:54:SER:CB	2.24	0.50
4:D:150:ASP:HB2	4:D:173:PRO:HG2	1.94	0.49
6:A:504:49M:H38	6:A:504:49M:H43	1.93	0.49
3:C:56:LYS:HG2	3:C:65:THR:HG22	1.94	0.49
1:E:153:ASP:OD2	6:E:504:49M:O2A	2.29	0.48
4:H:214:LEU:HD13	4:H:227:PRO:HG2	1.95	0.48
3:C:61:ARG:NH2	3:C:84:ASP:OD2	2.46	0.48
4:H:21:LEU:HD12	4:H:76:LEU:HD23	1.95	0.48
1:A:118:VAL:HG11	6:A:504:49M:H11	1.95	0.48
1:E:163:LEU:HB3	6:E:504:49M:H101	1.95	0.47
1:A:145:LEU:HB3	1:A:146:PRO:HD3	1.95	0.47
1:E:168:CYS:HB3	1:E:169:PRO:HD3	1.96	0.47
4:H:204:HIS:NE2	4:H:235:GLU:HG3	2.29	0.47
3:C:50:LEU:HD13	3:C:66:LEU:HB2	1.96	0.47
3:G:21:LEU:O	3:G:74:THR:HA	2.15	0.47
4:D:214:LEU:HD13	4:D:227:PRO:HG2	1.96	0.46
2:B:37:ILE:HG12	2:B:82:VAL:HG22	1.97	0.46
4:D:152:VAL:HG12	4:D:211:PHE:HA	1.96	0.46
1:E:216:LYS:N	1:E:217:PRO:HD2	2.31	0.45
3:G:8:PRO:O	3:G:108:THR:HG23	2.17	0.44
1:E:168:CYS:HB3	1:E:169:PRO:CD	2.47	0.44
4:D:46:ILE:HG22	4:D:47:HIS:HD2	1.83	0.44
4:D:46:ILE:HG22	4:D:47:HIS:CD2	2.52	0.44
4:H:36:ARG:HH21	4:H:87:SER:HB2	1.81	0.44
4:D:115:ARG:HA	4:D:222:GLN:HE22	1.83	0.44
1:E:39:ARG:HD3	2:F:53:ASP:OD2	2.18	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:H:78:LEU:HD23	4:H:85:GLN:OE1	2.18	0.43
3:G:13:VAL:HG22	3:G:110:LEU:HD11	2.01	0.43
1:E:114:SER:HB3	1:E:164:LEU:HD22	2.00	0.43
4:H:25:GLN:HE21	4:H:27:ASN:H	1.67	0.43
4:D:21:LEU:HD22	4:D:106:THR:HG21	2.00	0.43
4:D:144:ALA:HB2	4:D:209:VAL:HG21	2.00	0.43
3:G:50:LEU:HD13	3:G:66:LEU:HB2	2.01	0.43
3:G:202:PHE:HA	3:G:203:PRO:HD3	1.91	0.42
1:A:76:SER:HA	3:C:95:ARG:NH1	2.35	0.42
3:C:156:VAL:HG13	3:C:180:SER:HB2	2.01	0.42
4:H:132:ILE:HG23	4:H:195:ALA:HB1	2.02	0.42
4:H:127:PRO:HD2	4:H:198:TRP:CZ2	2.55	0.41
3:G:184:ASP:N	3:G:184:ASP:OD1	2.53	0.41
1:A:147:ILE:HA	1:A:147:ILE:HD13	1.92	0.41
1:A:207:VAL:HG22	1:A:250:THR:HG22	2.01	0.41
3:G:61:ARG:NH2	3:G:84:ASP:OD2	2.43	0.41
4:H:36:ARG:HB3	4:H:46:ILE:HD11	2.02	0.41
1:E:124:TYR:CZ	1:E:136:VAL:HG11	2.56	0.41
4:D:117:VAL:HG12	4:D:227:PRO:HB2	2.03	0.41
1:A:32:LEU:HD22	1:A:175:LEU:HD13	2.03	0.40
2:F:14:PRO:HA	2:F:15:PRO:HD3	1.85	0.40
3:C:128:ASP:HB3	3:C:131:SER:O	2.20	0.40
3:G:21:LEU:HD23	3:G:108:THR:HB	2.02	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	263/285 (92%)	257 (98%)	6 (2%)	0	100	100
1	E	263/285 (92%)	252 (96%)	10 (4%)	1 (0%)	39	80

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	94/99 (95%)	92 (98%)	1 (1%)	1 (1%)	17	62
2	F	94/99 (95%)	91 (97%)	2 (2%)	1 (1%)	17	62
3	C	199/209 (95%)	188 (94%)	11 (6%)	0	100	100
3	G	198/209 (95%)	188 (95%)	9 (4%)	1 (0%)	34	78
4	D	237/241 (98%)	235 (99%)	2 (1%)	0	100	100
4	H	237/241 (98%)	233 (98%)	4 (2%)	0	100	100
All	All	1585/1668 (95%)	1536 (97%)	45 (3%)	4 (0%)	46	85

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	F	47	PRO
1	E	56	GLY
2	B	47	PRO
3	G	27	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	218/249 (88%)	217 (100%)	1 (0%)	92	97
1	E	204/249 (82%)	204 (100%)	0	100	100
2	B	81/93 (87%)	81 (100%)	0	100	100
2	F	80/93 (86%)	79 (99%)	1 (1%)	76	92
3	C	174/188 (93%)	171 (98%)	3 (2%)	68	90
3	G	170/188 (90%)	169 (99%)	1 (1%)	90	97
4	D	198/208 (95%)	194 (98%)	4 (2%)	63	88
4	H	199/208 (96%)	195 (98%)	4 (2%)	63	88
All	All	1324/1476 (90%)	1310 (99%)	14 (1%)	82	94

All (14) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	170	LEU
3	C	141	ASP
3	C	181[A]	ASN
3	C	181[B]	ASN
4	D	25	GLN
4	D	55	THR
4	D	190	ARG
4	D	235	GLU
2	F	70	PHE
3	G	181	ASN
4	H	25	GLN
4	H	168	CYS
4	H	190	ARG
4	H	235	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (13) such sidechains are listed below:

Mol	Chain	Res	Type
3	C	31	ASN
3	C	125	GLN
3	C	193	ASN
4	D	47	HIS
4	D	116	ASN
4	D	217	ASN
4	D	222	GLN
1	E	36	GLN
1	E	117	HIS
3	G	31	ASN
3	G	117	GLN
4	H	25	GLN
4	H	222	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

8 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the chemical component dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
5	NAG	A	501	1	14,14,15	0.47	0	15,19,21	1.13	1 (6%)
5	NAG	A	502	1,5	14,14,15	0.46	0	15,19,21	0.94	1 (6%)
5	NAG	A	503	5	14,14,15	0.47	0	15,19,21	0.77	0
6	49M	A	504	-	62,63,63	0.30	0	64,72,72	0.72	2 (3%)
5	NAG	E	501	1	14,14,15	0.41	0	15,19,21	1.59	1 (6%)
5	NAG	E	502	1,5	14,14,15	0.50	0	15,19,21	0.72	0
5	NAG	E	503	5	14,14,15	0.53	0	15,19,21	1.01	1 (6%)
6	49M	E	504	-	62,63,63	0.29	0	64,72,72	0.91	3 (4%)

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	A	501	1	-	0/6/23/26	0/1/1/1
5	NAG	A	502	1,5	-	0/6/23/26	0/1/1/1
5	NAG	A	503	5	-	0/6/23/26	0/1/1/1
6	49M	A	504	-	-	0/61/81/81	0/1/1/1
5	NAG	E	501	1	-	0/6/23/26	0/1/1/1
5	NAG	E	502	1,5	-	0/6/23/26	0/1/1/1
5	NAG	E	503	5	-	0/6/23/26	0/1/1/1
6	49M	E	504	-	-	0/61/81/81	0/1/1/1

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	A	502	NAG	C1-O5-C5	2.27	115.12	112.25
6	A	504	49M	C5M-O6A-C1A	2.27	116.97	113.18
6	E	504	49M	C3A-C2A-C1A	2.37	114.02	109.32
6	A	504	49M	O6A-C1A-C2A	2.42	111.48	109.13
5	E	503	NAG	C4-C3-C2	2.45	115.04	111.23
6	E	504	49M	C5M-O6A-C1A	2.51	117.37	113.18
5	A	501	NAG	C1-O5-C5	3.62	116.84	112.25
6	E	504	49M	O6A-C1A-C2A	3.87	112.89	109.13
5	E	501	NAG	C1-O5-C5	5.33	119.01	112.25

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

3 monomers are involved in 9 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	504	49M	4	0
5	E	502	NAG	1	0
6	E	504	49M	4	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	267/285 (93%)	0.02	14 (5%) 31 18	64, 89, 139, 166	0
1	E	267/285 (93%)	0.16	24 (8%) 12 6	62, 97, 172, 189	0
2	B	96/99 (96%)	0.29	4 (4%) 40 26	79, 109, 124, 133	1 (1%)
2	F	96/99 (96%)	0.03	1 (1%) 84 75	68, 98, 125, 133	1 (1%)
3	C	202/209 (96%)	-0.13	1 (0%) 91 87	56, 77, 128, 141	0
3	G	202/209 (96%)	-0.25	0 100 100	59, 75, 122, 134	0
4	D	239/241 (99%)	-0.18	2 (0%) 87 80	61, 78, 108, 118	0
4	H	239/241 (99%)	-0.19	1 (0%) 93 90	65, 79, 110, 131	0
All	All	1608/1668 (96%)	-0.05	47 (2%) 55 41	56, 85, 135, 189	2 (0%)

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	E	256	GLY	5.7
1	A	262	ALA	5.4
1	E	259	ALA	4.1
1	E	253	VAL	4.0
1	E	222	TRP	4.0
1	E	260	GLY	3.8
1	A	222	TRP	3.6
4	D	179	ALA	3.5
4	H	240	ALA	3.5
1	E	108	PRO	3.4
1	E	255	ALA	3.3
1	E	276	ILE	3.3
1	E	231	GLY	3.2
1	E	275	ILE	3.1
1	E	251	LEU	3.1
1	E	252	ASP	3.1

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Mol	Chain	Res	Type	RSRZ
1	E	191	ALA	3.0
1	E	277	LEU	3.0
1	A	231	GLY	3.0
1	A	253	VAL	2.9
1	A	256	GLY	2.9
1	A	255	ALA	2.9
1	E	263	CYS	2.9
1	E	271	GLY	2.6
1	A	223	MET	2.5
1	E	278	TYR	2.5
2	B	18	GLY	2.5
1	A	251	LEU	2.4
1	E	262	ALA	2.4
1	E	221	MET	2.4
1	E	190	VAL	2.4
1	A	230	GLN	2.4
1	E	265	VAL	2.3
1	A	227	GLN	2.3
1	A	111	ALA	2.3
2	B	20	PRO	2.3
4	D	221	THR	2.3
1	E	258	GLU	2.3
2	B	82	VAL	2.3
1	E	230	GLN	2.3
3	C	191	PHE	2.2
1	A	229	GLN	2.2
2	B	78	TYR	2.1
1	E	107	TYR	2.1
1	A	250	THR	2.0
2	F	71	THR	2.0
1	A	233	HIS	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 carbohydrates 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. 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	49M	E	504	63/63	0.91	0.36	4.23	71,79,94,98	0
6	49M	A	504	63/63	0.93	0.30	2.65	62,67,72,73	0
5	NAG	E	501	14/15	0.90	0.24	0.55	89,91,93,94	0
5	NAG	A	501	14/15	0.89	0.23	-	87,89,91,92	0
5	NAG	E	502	14/15	0.92	0.16	-	81,83,86,89	0
5	NAG	A	502	14/15	0.91	0.21	-	74,76,78,79	0
5	NAG	A	503	14/15	0.94	0.20	-	82,85,87,87	0
5	NAG	E	503	14/15	0.90	0.29	-	94,100,103,105	0

6.5 Other polymers [i](#)

There are no such residues in this entry.