



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 1, 2016 – 02:25 AM GMT

PDB ID : 2H0B
Title : Crystal Structure of the second LNS/LG domain from Neurexin 1 alpha
Authors : Sheckler, L.R.; Henry, L.; Sugita, S.; Sudhof, T.C.; Rudenko, G.
Deposited on : 2006-05-14
Resolution : 2.10 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

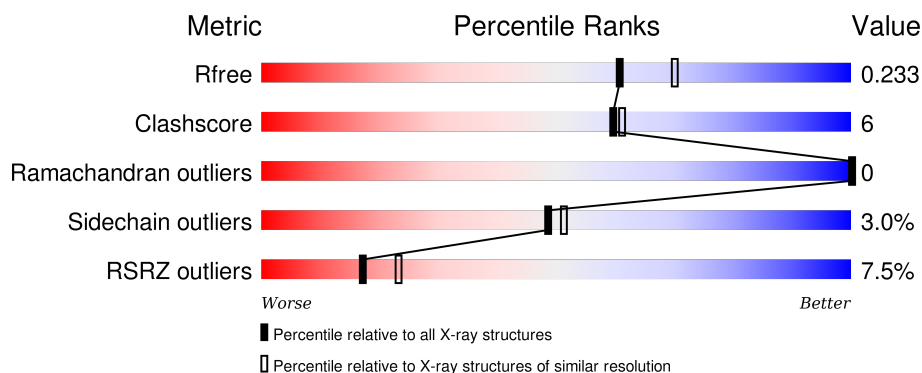
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.10 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	3939 (2.10-2.10)
Clashscore	102246	4460 (2.10-2.10)
Ramachandran outliers	100387	4413 (2.10-2.10)
Sidechain outliers	100360	4414 (2.10-2.10)
RSRZ outliers	91569	3948 (2.10-2.10)

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	184	<div> <div>7%</div> <div> <div></div> <div>84%</div> <div>13%</div> <div>...</div> </div> </div>
1	B	184	<div> <div>5%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>...</div> </div> </div>
1	C	184	<div> <div>11%</div> <div> <div></div> <div>85%</div> <div>13%</div> <div>...</div> </div> </div>
1	D	184	<div> <div>6%</div> <div> <div></div> <div>83%</div> <div>14%</div> <div>...</div> </div> </div>

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
3	GOL	A	4503	-	-	-	X
3	GOL	D	4501	-	-	-	X

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6025 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 Neurexin-1-alpha.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	181	Total	C	N	O	S	Se	0	0	0
			1390	875	233	276	2	4			
1	B	181	Total	C	N	O	S	Se	0	0	0
			1387	872	232	277	2	4			
1	C	183	Total	C	N	O	S	Se	0	0	0
			1385	870	232	277	2	4			
1	D	182	Total	C	N	O	S	Se	0	0	0
			1391	871	234	280	2	4			

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	277	GLY	-	CLONING ARTIFACT	UNP Q28146
A	278	SER	-	CLONING ARTIFACT	UNP Q28146
A	321	MSE	MET	MODIFIED RESIDUE	UNP Q28146
A	414	MSE	MET	MODIFIED RESIDUE	UNP Q28146
A	442	MSE	MET	MODIFIED RESIDUE	UNP Q28146
A	470	MSE	MET	MODIFIED RESIDUE	UNP Q28146
B	277	GLY	-	CLONING ARTIFACT	UNP Q28146
B	278	SER	-	CLONING ARTIFACT	UNP Q28146
B	321	MSE	MET	MODIFIED RESIDUE	UNP Q28146
B	414	MSE	MET	MODIFIED RESIDUE	UNP Q28146
B	442	MSE	MET	MODIFIED RESIDUE	UNP Q28146
B	470	MSE	MET	MODIFIED RESIDUE	UNP Q28146
C	277	GLY	-	CLONING ARTIFACT	UNP Q28146
C	278	SER	-	CLONING ARTIFACT	UNP Q28146
C	321	MSE	MET	MODIFIED RESIDUE	UNP Q28146
C	414	MSE	MET	MODIFIED RESIDUE	UNP Q28146
C	442	MSE	MET	MODIFIED RESIDUE	UNP Q28146
C	470	MSE	MET	MODIFIED RESIDUE	UNP Q28146
D	277	GLY	-	CLONING ARTIFACT	UNP Q28146
D	278	SER	-	CLONING ARTIFACT	UNP Q28146
D	321	MSE	MET	MODIFIED RESIDUE	UNP Q28146

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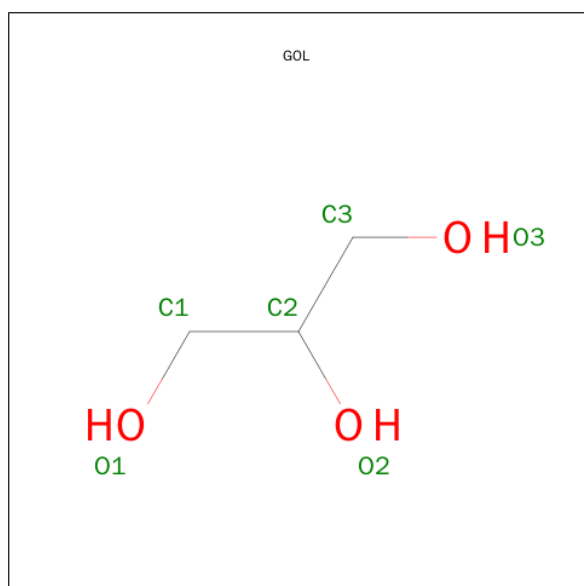
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Chain	Residue	Modelled	Actual	Comment	Reference
D	414	MSE	MET	MODIFIED RESIDUE	UNP Q28146
D	442	MSE	MET	MODIFIED RESIDUE	UNP Q28146
D	470	MSE	MET	MODIFIED RESIDUE	UNP Q28146

- Molecule 2 is CALCIUM ION (three-letter code: CA) (formula: Ca).

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

- Molecule 3 is GLYCEROL (three-letter code: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C O 6 3 3	0	0
3	D	1	Total C O 6 3 3	0	0
3	B	1	Total C O 6 3 3	0	0
3	A	1	Total C O 6 3 3	0	0

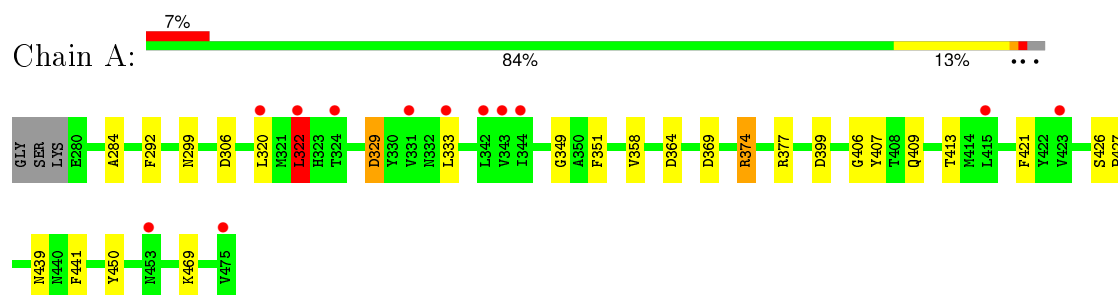
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	115	Total 115	O 115	0	0
4	B	130	Total 130	O 130	0	0
4	C	90	Total 90	O 90	0	0
4	D	109	Total 109	O 109	0	0

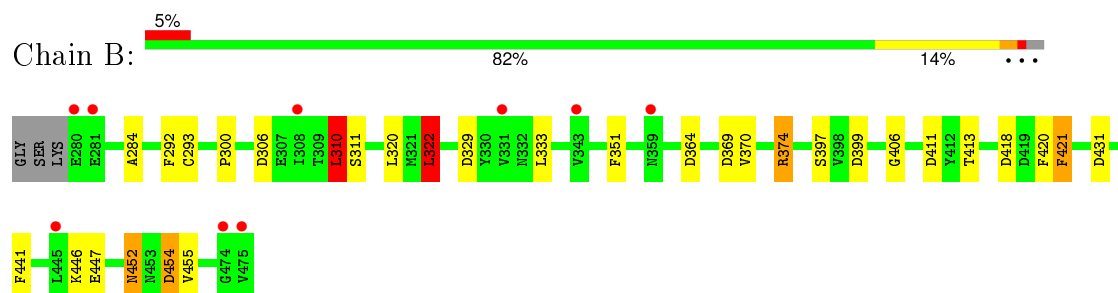
3 Residue-property plots [i](#)

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 ($\text{RSRZ} > 2$). 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.

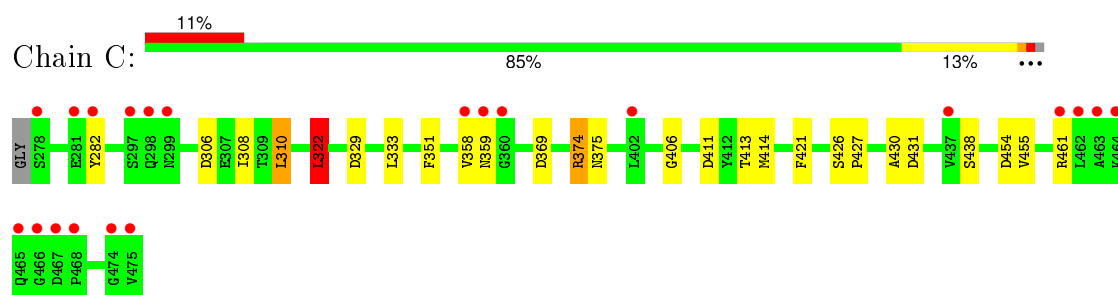
• Molecule 1: Neurexin-1-alpha



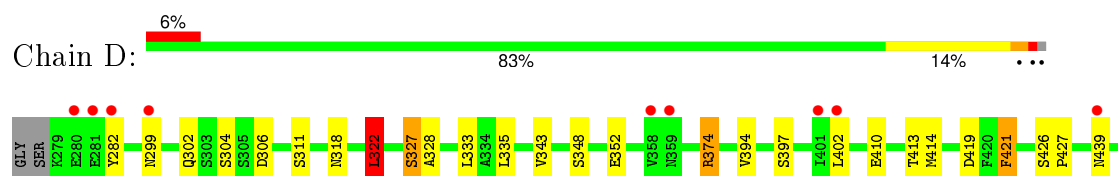
• Molecule 1: Neurexin-1-alpha

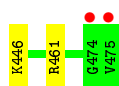


• Molecule 1: Neurexin-1-alpha



• Molecule 1: Neurexin-1-alpha





4 Data and refinement statistics

Property	Value	Source
Space group	P 1	Depositor
Cell constants a, b, c, α , β , γ	60.03Å 61.59Å 66.64Å 78.47° 78.74° 84.68°	Depositor
Resolution (Å)	20.00 – 2.10 34.53 – 2.10	Depositor EDS
% Data completeness (in resolution range)	100.0 (20.00-2.10) 91.3 (34.53-2.10)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.31 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.2.0003	Depositor
R, R_{free}	0.184 , 0.233 0.185 , 0.233	Depositor DCC
R_{free} test set	2577 reflections (5.27%)	DCC
Wilson B-factor (Å ²)	26.3	Xtriage
Anisotropy	0.399	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 41.6	EDS
Estimated twinning fraction	0.014 for -k,-h,-l	Xtriage
L-test for twinning ²	$\langle L \rangle = 0.50$, $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 51540 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	6025	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.23% 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: GOL, CA

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	0.54	0/1412	0.87	6/1906 (0.3%)
1	B	0.58	0/1409	0.90	10/1902 (0.5%)
1	C	0.47	0/1407	0.81	6/1902 (0.3%)
1	D	0.50	0/1412	0.84	3/1904 (0.2%)
All	All	0.52	0/5640	0.86	25/7614 (0.3%)

There are no bond length outliers.

All (25) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	322	LEU	CA-CB-CG	7.70	133.02	115.30
1	B	310	LEU	CA-CB-CG	7.37	132.26	115.30
1	B	374	ARG	NE-CZ-NH2	-7.25	116.67	120.30
1	D	322	LEU	CA-CB-CG	7.20	131.85	115.30
1	D	374	ARG	NE-CZ-NH2	-7.11	116.74	120.30
1	A	322	LEU	CA-CB-CG	6.98	131.36	115.30
1	A	374	ARG	NE-CZ-NH2	-6.98	116.81	120.30
1	B	322	LEU	CA-CB-CG	6.32	129.84	115.30
1	A	369	ASP	CB-CG-OD2	6.08	123.77	118.30
1	C	329	ASP	CB-CG-OD2	6.08	123.77	118.30
1	C	310	LEU	CA-CB-CG	5.96	129.00	115.30
1	C	454	ASP	CB-CG-OD2	5.81	123.53	118.30
1	B	454	ASP	CB-CG-OD2	5.80	123.52	118.30
1	A	364	ASP	CB-CG-OD2	5.71	123.44	118.30
1	B	364	ASP	CB-CG-OD2	5.69	123.42	118.30
1	B	411	ASP	CB-CG-OD2	5.61	123.35	118.30
1	C	431	ASP	CB-CG-OD2	5.55	123.29	118.30
1	A	329	ASP	CB-CG-OD2	5.43	123.19	118.30
1	B	374	ARG	NE-CZ-NH1	5.42	123.01	120.30
1	D	419	ASP	CB-CG-OD2	5.34	123.11	118.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	369	ASP	CB-CG-OD2	5.34	123.10	118.30
1	B	431	ASP	CB-CG-OD2	5.17	122.95	118.30
1	C	369	ASP	CB-CG-OD2	5.12	122.91	118.30
1	A	399	ASP	CB-CG-OD2	5.12	122.91	118.30
1	B	399	ASP	CB-CG-OD2	5.07	122.86	118.30

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	1390	0	1352	16	0
1	B	1387	0	1341	17	0
1	C	1385	0	1320	12	0
1	D	1391	0	1343	18	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	12	0	16	1	0
3	B	6	0	8	0	0
3	D	6	0	8	0	0
4	A	115	0	0	1	0
4	B	130	0	0	0	0
4	C	90	0	0	1	0
4	D	109	0	0	4	0
All	All	6025	0	5388	62	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 6.

All (62) 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:D:304:SER:HB3	1:D:414:MSE:HE1	1.53	0.90
1:B:452:ASN:ND2	1:B:454:ASP:H	1.84	0.76
1:D:374:ARG:NH2	1:D:413:THR:O	2.22	0.73
1:B:452:ASN:HD22	1:B:454:ASP:H	1.38	0.67
1:D:304:SER:HB3	1:D:414:MSE:CE	2.24	0.66
1:A:322:LEU:HD22	1:A:333:LEU:HB3	1.78	0.66
1:C:322:LEU:HD22	1:C:333:LEU:HB3	1.78	0.65
1:C:374:ARG:NH2	1:C:413:THR:O	2.30	0.64
1:A:299:ASN:HD21	1:A:469:LYS:HZ1	1.47	0.62
1:B:452:ASN:HD22	1:B:452:ASN:C	2.04	0.61
1:A:439:ASN:ND2	4:A:5035:HOH:O	2.34	0.60
1:B:322:LEU:HD22	1:B:333:LEU:HB3	1.84	0.59
1:B:374:ARG:NH2	1:B:413:THR:O	2.35	0.59
1:D:302:GLN:HG3	4:D:5423:HOH:O	2.02	0.57
1:A:299:ASN:ND2	1:A:469:LYS:NZ	2.51	0.57
1:A:306:ASP:OD2	1:A:374:ARG:NH1	2.35	0.56
1:A:299:ASN:HD21	1:A:469:LYS:NZ	2.03	0.56
1:B:300:PRO:HB3	1:B:418:ASP:HB3	1.87	0.56
1:C:351:PHE:CE1	1:C:406:GLY:HA3	2.42	0.55
1:D:322:LEU:HD22	1:D:333:LEU:HB3	1.89	0.54
1:B:310:LEU:HD22	1:B:370:VAL:HB	1.89	0.53
1:A:374:ARG:NH2	1:A:413:THR:O	2.41	0.53
1:C:282:TYR:CZ	1:C:461:ARG:HG3	2.44	0.52
1:C:426:SER:HB2	1:C:427:PRO:HD2	1.92	0.52
1:B:306:ASP:OD2	1:B:374:ARG:NH1	2.41	0.52
1:D:439:ASN:ND2	4:D:5301:HOH:O	2.43	0.51
1:A:351:PHE:CZ	1:A:406:GLY:HA3	2.49	0.48
1:D:348:SER:HB2	1:D:410:GLU:HB2	1.95	0.48
1:B:322:LEU:HD23	1:B:322:LEU:C	2.33	0.48
1:D:282:TYR:CZ	1:D:461:ARG:HG3	2.48	0.48
1:A:299:ASN:ND2	1:A:469:LYS:HZ3	2.13	0.46
1:D:426:SER:HB2	1:D:427:PRO:HD2	1.97	0.46
1:A:377:ARG:O	1:A:407:TYR:HA	2.16	0.46
1:D:397:SER:HA	1:D:402:LEU:O	2.14	0.46
1:C:358:VAL:HG12	1:C:359:ASN:HD22	1.80	0.46
1:B:452:ASN:ND2	1:B:455:VAL:H	2.14	0.45
1:B:351:PHE:CZ	1:B:406:GLY:HA3	2.51	0.45
1:A:284:ALA:CB	1:A:292:PHE:HZ	2.30	0.45
1:D:343:VAL:HG22	1:D:352:GLU:HG2	1.98	0.44
1:D:311:SER:HB2	1:D:446:LYS:HG2	1.99	0.44
1:C:308:ILE:HG21	1:C:322:LEU:HD11	1.99	0.44
1:A:426:SER:HB2	1:A:427:PRO:HD2	1.99	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:414:MSE:HE1	4:D:5329:HOH:O	2.18	0.43
1:A:306:ASP:OD1	1:A:450:TYR:OH	2.25	0.43
1:C:375:ASN:OD1	1:D:439:ASN:ND2	2.52	0.43
1:B:320:LEU:O	1:B:441:PHE:HB2	2.19	0.42
1:B:446:LYS:HG3	1:B:447:GLU:HG2	2.00	0.42
1:D:327:SER:HB3	1:D:328:ALA:H	1.58	0.42
1:D:322:LEU:HG	1:D:421:PHE:HE2	1.84	0.42
1:B:293:CYS:SG	1:B:420:PHE:CD2	3.13	0.42
1:B:284:ALA:CB	1:B:292:PHE:HZ	2.33	0.41
1:A:349:GLY:HA3	1:A:409:GLN:HB2	2.02	0.41
1:A:320:LEU:O	1:A:441:PHE:HB2	2.20	0.41
1:D:299:ASN:HB3	4:D:5311:HOH:O	2.19	0.41
1:A:358:VAL:H	3:A:4503:GOL:C1	2.33	0.41
1:B:322:LEU:HG	1:B:421:PHE:HE2	1.85	0.41
1:C:411:ASP:HB2	4:C:5130:HOH:O	2.20	0.41
1:D:374:ARG:HG3	1:D:394:VAL:HG22	2.03	0.41
1:C:430:ALA:HB2	1:C:438:SER:HB2	2.01	0.40
1:B:311:SER:HB2	1:B:446:LYS:HG2	2.04	0.40
1:C:306:ASP:OD2	1:C:374:ARG:NH1	2.50	0.40
1:D:318:ASN:HA	1:D:335:LEU:O	2.21	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	179/184 (97%)	171 (96%)	8 (4%)	0	100	100
1	B	179/184 (97%)	170 (95%)	9 (5%)	0	100	100
1	C	181/184 (98%)	172 (95%)	9 (5%)	0	100	100
1	D	180/184 (98%)	173 (96%)	7 (4%)	0	100	100
All	All	719/736 (98%)	686 (95%)	33 (5%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	154/154 (100%)	151 (98%)	3 (2%)	65	70
1	B	153/154 (99%)	147 (96%)	6 (4%)	39	39
1	C	150/154 (97%)	145 (97%)	5 (3%)	45	47
1	D	153/154 (99%)	149 (97%)	4 (3%)	54	58
All	All	610/616 (99%)	592 (97%)	18 (3%)	48	51

All (18) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	322	LEU
1	A	329	ASP
1	A	421	PHE
1	B	310	LEU
1	B	322	LEU
1	B	329	ASP
1	B	397	SER
1	B	421	PHE
1	B	452	ASN
1	C	310	LEU
1	C	322	LEU
1	C	374	ARG
1	C	421	PHE
1	C	455	VAL
1	D	306	ASP
1	D	322	LEU
1	D	327	SER
1	D	421	PHE

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

Mol	Chain	Res	Type
1	A	299	ASN
1	A	302	GLN
1	A	440	ASN
1	B	332	ASN
1	B	365	ASN
1	B	375	ASN
1	B	440	ASN
1	B	452	ASN
1	C	332	ASN
1	C	359	ASN
1	C	440	ASN
1	D	332	ASN
1	D	440	ASN

5.3.3 RNA [i](#)

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 carbohydrates in this entry.

5.6 Ligand geometry [i](#)

Of 8 ligands modelled in this entry, 4 are monoatomic - leaving 4 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
3	GOL	A	4500	-	5,5,5	0.36	0	5,5,5	0.60	0
3	GOL	A	4503	-	5,5,5	0.21	0	5,5,5	0.45	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	B	4502	-	5,5,5	0.30	0	5,5,5	0.65	0
3	GOL	D	4501	-	5,5,5	0.34	0	5,5,5	0.99	0

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the chemical component dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	GOL	A	4500	-	-	0/4/4/4	0/0/0/0
3	GOL	A	4503	-	-	0/4/4/4	0/0/0/0
3	GOL	B	4502	-	-	0/4/4/4	0/0/0/0
3	GOL	D	4501	-	-	0/4/4/4	0/0/0/0

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.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	4503	GOL	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	177/184 (96%)	0.53	12 (6%) 20 28	31, 35, 40, 47	0
1	B	177/184 (96%)	0.46	9 (5%) 32 40	30, 35, 41, 51	0
1	C	179/184 (97%)	0.77	21 (11%) 6 8	28, 36, 42, 50	0
1	D	178/184 (96%)	0.38	11 (6%) 24 32	25, 35, 43, 55	0
All	All	711/736 (96%)	0.53	53 (7%) 17 23	25, 36, 41, 55	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	359	ASN	6.6
1	D	359	ASN	6.3
1	D	475	VAL	5.6
1	C	475	VAL	5.4
1	C	297	SER	4.8
1	C	468	PRO	4.2
1	B	475	VAL	4.0
1	D	358	VAL	3.9
1	A	331	VAL	3.8
1	C	299	ASN	3.8
1	A	453	ASN	3.5
1	C	358	VAL	3.5
1	C	467	ASP	3.4
1	B	280	GLU	3.3
1	C	298	GLN	3.3
1	C	281	GLU	3.2
1	D	401	ILE	3.2
1	B	359	ASN	2.9
1	A	344	ILE	2.8
1	D	474	GLY	2.7
1	A	475	VAL	2.7

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Mol	Chain	Res	Type	RSRZ
1	C	466	GLY	2.7
1	B	331	VAL	2.7
1	D	280	GLU	2.7
1	D	282	TYR	2.6
1	D	402	LEU	2.6
1	C	278	SER	2.6
1	A	333	LEU	2.6
1	C	464	LYS	2.5
1	C	282	TYR	2.5
1	B	474	GLY	2.5
1	C	474	GLY	2.5
1	A	423	VAL	2.5
1	D	299	ASN	2.4
1	C	463	ALA	2.4
1	C	462	LEU	2.4
1	B	281	GLU	2.4
1	A	342	LEU	2.3
1	C	437	VAL	2.3
1	C	360	GLY	2.3
1	A	324	THR	2.3
1	D	439	ASN	2.3
1	A	320	LEU	2.3
1	C	402	LEU	2.2
1	D	281	GLU	2.1
1	A	343	VAL	2.1
1	C	461	ARG	2.1
1	A	322	LEU	2.1
1	A	415	LEU	2.1
1	B	343	VAL	2.1
1	B	445	LEU	2.0
1	C	465	GLN	2.0
1	B	308	ILE	2.0

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

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

6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

6.4 Ligands

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
3	GOL	A	4503	6/6	0.81	0.25	4.52	57,57,58,59	0
3	GOL	D	4501	6/6	0.85	0.25	2.56	32,36,38,38	0
3	GOL	B	4502	6/6	0.93	0.14	-0.19	33,36,37,40	0
3	GOL	A	4500	6/6	0.94	0.11	-0.77	32,34,36,40	0
2	CA	C	3000	1/1	1.00	0.06	-2.64	25,25,25,25	0
2	CA	A	1000	1/1	0.99	0.04	-2.66	27,27,27,27	0
2	CA	B	2000	1/1	0.98	0.03	-3.82	24,24,24,24	0
2	CA	D	4000	1/1	0.99	0.06	-4.26	27,27,27,27	0

6.5 Other polymers

There are no such residues in this entry.