



Full wwPDB X-ray Structure Validation Report ⓘ

Feb 6, 2017 – 01:34 PM EST

PDB ID : 5WPZ
Title : Crystal structure of MNDA PYD with MBP tag
Authors : Jin, T.C.; Xiao, T.S.
Deposited on : 2016-11-22
Resolution : 2.00 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 : 1.7.1 (RC1), CSD as537be (2016)
Xtriage (Phenix) : 1.9-1692
EDS : rb-20028442
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) : rb-20028442

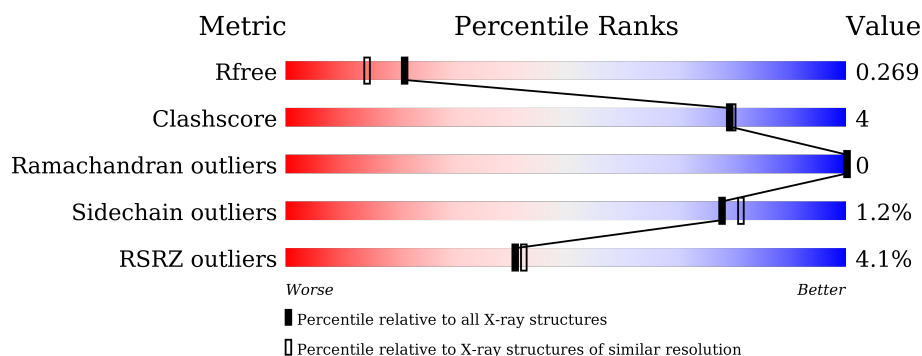
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.00 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	6249 (2.00-2.00)
Clashscore	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)
RSRZ outliers	91569	6262 (2.00-2.00)

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	477	<div> <div>0%</div> <div>88%</div> <div>8%</div> <div>.</div> </div>
1	B	477	<div> <div>2%</div> <div>89%</div> <div>7%</div> <div>..</div> </div>
1	C	477	<div> <div>3%</div> <div>89%</div> <div>8%</div> <div>.</div> </div>
1	D	477	<div> <div>3%</div> <div>86%</div> <div>10%</div> <div>.</div> </div>
1	E	477	<div> <div>5%</div> <div>87%</div> <div>9%</div> <div>.</div> </div>
1	F	477	<div> <div>9%</div> <div>84%</div> <div>11%</div> <div>.</div> </div>

2 Entry composition [i](#)

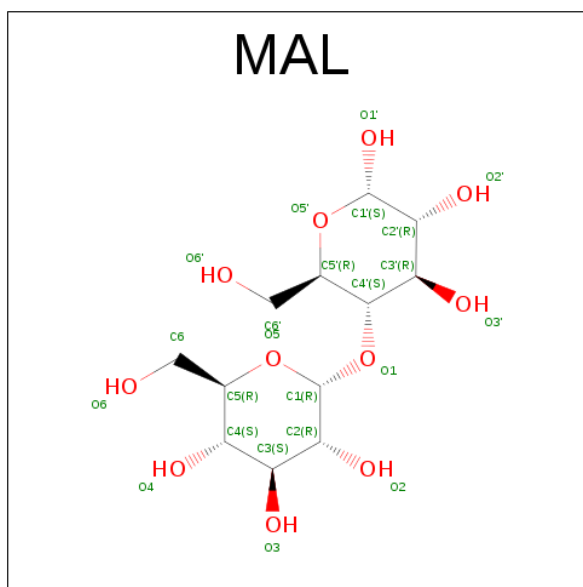
There are 3 unique types of molecules in this entry. The entry contains 22797 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 MBP-hMNDA-PYD.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	460	Total	C	N	O	S	0	1	0
			3576	2304	585	676	11			
1	B	462	Total	C	N	O	S	0	0	0
			3585	2311	587	676	11			
1	C	461	Total	C	N	O	S	0	3	0
			3599	2319	591	678	11			
1	D	459	Total	C	N	O	S	0	2	0
			3579	2306	587	675	11			
1	E	463	Total	C	N	O	S	0	1	0
			3599	2319	590	679	11			
1	F	459	Total	C	N	O	S	0	2	0
			3579	2306	587	675	11			

- Molecule 2 is MALTOSE (three-letter code: MAL) (formula: $C_{12}H_{22}O_{11}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			23	12	11		
2	B	1	Total	C	O	0	0
			23	12	11		
2	C	1	Total	C	O	0	0
			23	12	11		
2	D	1	Total	C	O	0	0
			23	12	11		
2	E	1	Total	C	O	0	0
			23	12	11		
2	F	1	Total	C	O	0	0
			23	12	11		

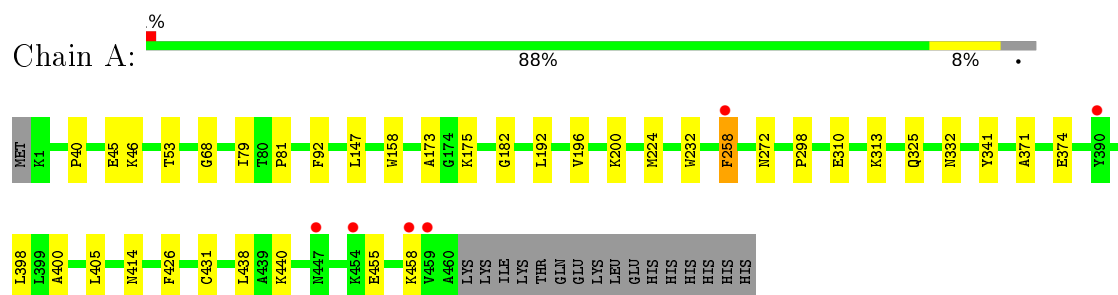
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	240	Total	O	0	0
			240	240		
3	B	239	Total	O	0	0
			239	239		
3	C	189	Total	O	0	0
			189	189		
3	D	173	Total	O	0	0
			173	173		
3	E	159	Total	O	0	0
			159	159		
3	F	142	Total	O	0	0
			142	142		

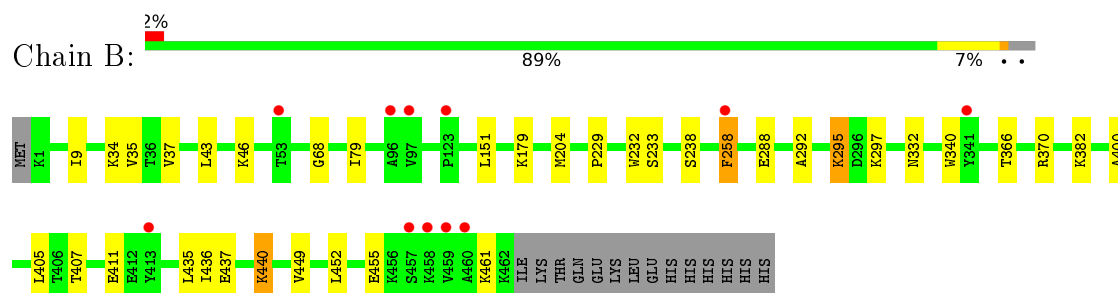
3 Residue-property plots [i](#)

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.

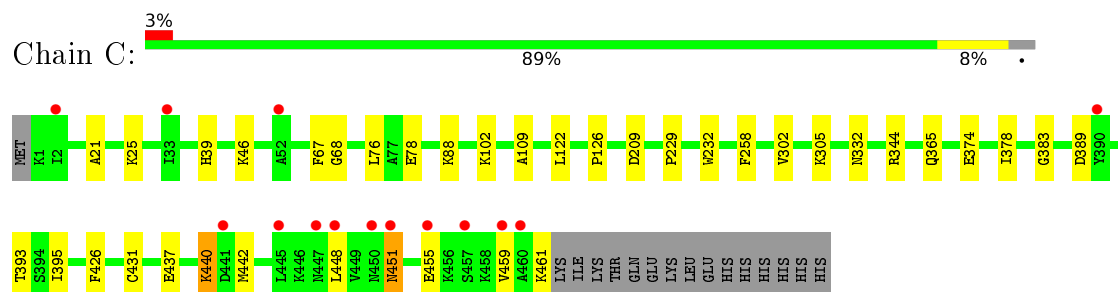
• Molecule 1: MBP-hMNDA-PYD



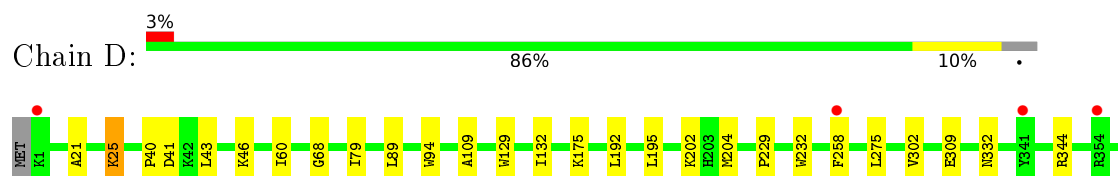
• Molecule 1: MBP-hMNDA-PYD



• Molecule 1: MBP-hMNDA-PYD

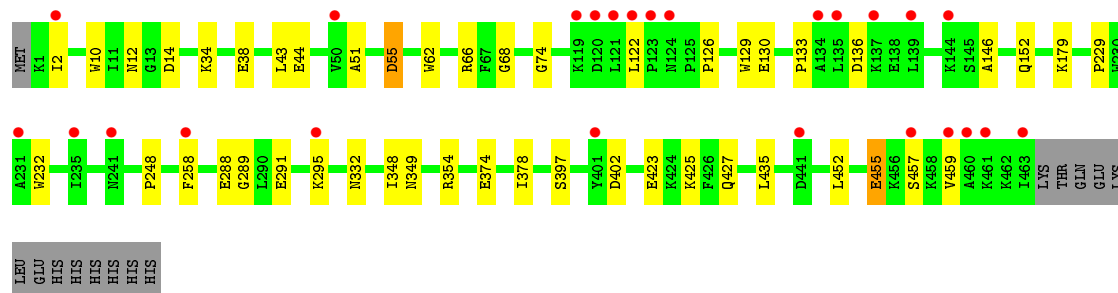
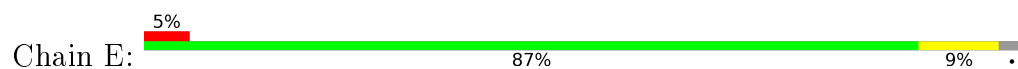


• Molecule 1: MBP-hMNDA-PYD

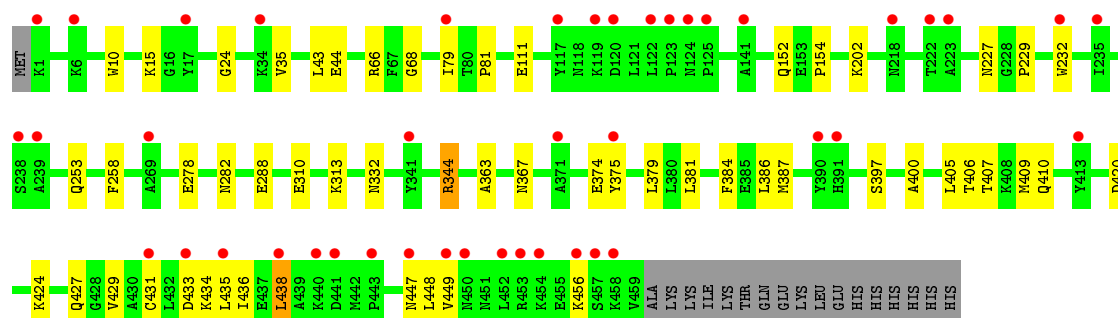
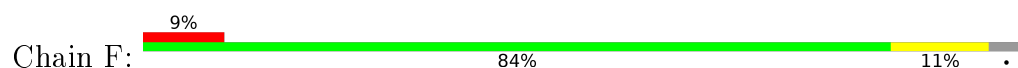




• Molecule 1: MBP-hMNDA-PYD



• Molecule 1: MBP-hMNDA-PYD



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 2	Depositor
Cell constants a, b, c, α , β , γ	192.15Å 235.61Å 73.76Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.36 – 2.00 48.36 – 2.00	Depositor EDS
% Data completeness (in resolution range)	89.0 (48.36-2.00) 89.0 (48.36-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.91 (at 2.00Å)	Xtriage
Refinement program	PHENIX (dev_2481: ???)	Depositor
R, R_{free}	0.221 , 0.269 0.219 , 0.269	Depositor DCC
R_{free} test set	9994 reflections (4.96%)	DCC
Wilson B-factor (Å ²)	30.3	Xtriage
Anisotropy	0.853	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 38.5	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.93	EDS
Total number of atoms	22797	wwPDB-VP
Average B, all atoms (Å ²)	37.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 7.75% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: MAL

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	0.26	0/3654	0.43	0/4950
1	B	0.27	0/3663	0.43	0/4960
1	C	0.27	0/3683	0.43	0/4988
1	D	0.26	0/3660	0.43	0/4957
1	E	0.26	0/3677	0.43	0/4979
1	F	0.27	0/3660	0.43	0/4957
All	All	0.26	0/21997	0.43	0/29791

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	3576	0	3597	27	0
1	B	3585	0	3616	21	0
1	C	3599	0	3631	23	0
1	D	3579	0	3605	26	0
1	E	3599	0	3625	27	0
1	F	3579	0	3605	32	0
2	A	23	0	22	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	B	23	0	22	0	0
2	C	23	0	22	1	0
2	D	23	0	22	0	0
2	E	23	0	22	0	0
2	F	23	0	22	0	0
3	A	240	0	0	7	0
3	B	239	0	0	5	0
3	C	189	0	0	2	0
3	D	173	0	0	1	0
3	E	159	0	0	1	0
3	F	142	0	0	3	0
All	All	22797	0	21811	154	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 4.

All (154) 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:C:437:GLU:O	1:C:440:LYS:HE3	1.77	0.83
1:F:436:ILE:HG12	1:F:449:VAL:HG13	1.67	0.74
1:B:295:LYS:NZ	3:B:601:HOH:O	2.20	0.72
1:D:399:LEU:HB3	1:D:403:LEU:HD12	1.71	0.72
1:F:407:THR:HA	1:F:410:GLN:HB2	1.72	0.71
1:F:79:ILE:HD12	1:F:81:PRO:HD3	1.73	0.71
1:F:24:GLY:HA3	1:F:35:VAL:HG21	1.77	0.66
1:A:173:ALA:O	3:A:601:HOH:O	2.15	0.64
1:F:44:GLU:HB2	1:F:66:ARG:HD3	1.81	0.62
1:A:440:LYS:NZ	3:A:608:HOH:O	2.33	0.62
1:E:374:GLU:CD	1:E:374:GLU:H	2.03	0.62
1:F:400:ALA:HA	1:F:405:LEU:HB2	1.82	0.61
1:B:382:LYS:HG3	1:B:455:GLU:HG2	1.82	0.61
1:D:383:GLY:HA2	1:D:448:LEU:HD21	1.83	0.60
1:C:378:ILE:HD11	1:C:459:VAL:HG11	1.84	0.60
1:B:297:LYS:NZ	3:B:608:HOH:O	2.36	0.59
1:F:447:ASN:OD1	1:F:448:LEU:N	2.35	0.59
1:B:400:ALA:HA	1:B:405:LEU:HB2	1.85	0.58
1:B:435:LEU:HD23	1:B:452:LEU:HD11	1.85	0.58
1:A:147:LEU:HB2	1:A:224:MET:HE3	1.86	0.58
1:C:67:PHE:HD1	1:C:76:LEU:HD11	1.68	0.58
1:E:291:GLU:O	1:E:295:LYS:HB2	2.04	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:365:GLN:NE2	3:C:606:HOH:O	2.31	0.57
1:D:400:ALA:HA	1:D:405:LEU:HB2	1.86	0.57
1:D:21:ALA:O	1:D:25:LYS:HD2	2.04	0.57
1:E:10:TRP:CD1	1:E:38:GLU:HG3	2.40	0.57
1:A:325:GLN:NE2	3:A:610:HOH:O	2.36	0.57
1:F:313:LYS:NZ	3:F:604:HOH:O	2.27	0.57
1:F:253:GLN:NE2	3:F:610:HOH:O	2.38	0.57
1:F:154:PRO:HD3	1:F:344[B]:ARG:HG3	1.87	0.56
1:A:310:GLU:OE1	1:A:313:LYS:NZ	2.30	0.56
1:C:21:ALA:O	1:C:25:LYS:HG2	2.06	0.56
1:A:398:LEU:HD23	1:A:438:LEU:HD22	1.87	0.55
1:A:400:ALA:HA	1:A:405:LEU:HB2	1.89	0.55
1:D:41:ASP:O	1:D:46:LYS:NZ	2.28	0.54
1:F:10:TRP:HB3	1:F:43:LEU:HD11	1.89	0.54
1:D:455:GLU:OE2	1:D:458:LYS:HE3	2.08	0.53
1:C:378:ILE:HG23	1:C:455:GLU:HG2	1.91	0.53
1:E:68:GLY:HA3	1:E:332:ASN:O	2.08	0.53
1:D:436:ILE:HG23	1:D:449:VAL:HG13	1.91	0.53
1:D:68:GLY:HA3	1:D:332:ASN:O	2.07	0.53
1:D:451:ASN:HA	1:D:454:LYS:HG2	1.90	0.53
1:F:44:GLU:OE1	1:F:44:GLU:N	2.42	0.53
1:F:363:ALA:O	1:F:367:ASN:HB2	2.08	0.52
1:D:450:ASN:OD1	1:D:453:ARG:NH1	2.43	0.52
1:E:10:TRP:HB3	1:E:43:LEU:HD11	1.91	0.52
1:B:436:ILE:HG23	1:B:449:VAL:HG13	1.90	0.52
1:D:202:LYS:NZ	3:D:610:HOH:O	2.42	0.52
1:E:44:GLU:N	1:E:44:GLU:OE1	2.34	0.52
1:D:388:ASP:OD1	1:D:391:HIS:ND1	2.44	0.51
1:E:74:GLY:O	3:E:601:HOH:O	2.19	0.51
1:C:395:ILE:HD11	1:C:442:MET:HE2	1.91	0.51
1:C:383:GLY:HA2	1:C:448:LEU:HD21	1.94	0.50
1:E:402:ASP:O	1:E:425:LYS:NZ	2.42	0.50
1:A:68:GLY:HA3	1:A:332:ASN:O	2.11	0.50
1:A:45:GLU:OE2	1:A:341:TYR:OH	2.27	0.50
1:F:374:GLU:N	1:F:374:GLU:OE1	2.39	0.50
1:B:34:LYS:HD2	1:B:35:VAL:N	2.27	0.49
1:B:407:THR:O	1:B:411:GLU:HG3	2.12	0.49
1:C:451:ASN:O	1:C:455:GLU:HB2	2.13	0.49
1:B:68:GLY:HA3	1:B:332:ASN:O	2.11	0.49
1:B:179:LYS:NZ	3:B:603:HOH:O	2.35	0.49
1:E:374:GLU:N	1:E:374:GLU:OE1	2.34	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:232:TRP:HB2	1:A:298:PRO:HG2	1.95	0.48
1:D:455:GLU:OE2	1:D:455:GLU:HA	2.13	0.48
1:F:310:GLU:HA	1:F:313:LYS:HE3	1.95	0.48
1:F:434:LYS:O	1:F:438:LEU:HG	2.14	0.48
1:B:9:ILE:HB	1:B:37:VAL:HG22	1.96	0.48
1:D:129:TRP:HA	1:D:132:ILE:HD13	1.96	0.48
1:E:288:GLU:HG2	1:E:289:GLY:N	2.29	0.48
1:B:79:ILE:O	3:B:602:HOH:O	2.20	0.48
1:F:420:ASP:O	1:F:424:LYS:HG3	2.14	0.48
1:E:374:GLU:O	1:E:378:ILE:HD12	2.14	0.48
1:A:192:LEU:O	1:A:196:VAL:HG23	2.14	0.47
1:C:344[B]:ARG:NH2	2:C:501:MAL:H62	2.29	0.47
1:C:455:GLU:O	1:C:459:VAL:HG22	2.14	0.47
1:D:195:LEU:HD12	1:D:204:MET:HE1	1.97	0.47
1:F:288:GLU:CD	1:F:288:GLU:H	2.16	0.47
1:F:278:GLU:OE2	1:F:282:ASN:ND2	2.38	0.47
1:F:381:LEU:O	1:F:386:LEU:HD13	2.13	0.47
1:C:39:HIS:ND1	1:C:39:HIS:O	2.48	0.46
1:D:376:LYS:HG3	1:D:423:GLU:HG3	1.98	0.46
1:D:43:LEU:HD13	1:D:60:ILE:HD11	1.97	0.46
1:D:275:LEU:H	1:D:275:LEU:HD12	1.81	0.46
1:A:40:PRO:HB2	1:A:46:LYS:HE3	1.97	0.46
1:A:272:ASN:OD1	3:A:602:HOH:O	2.20	0.46
1:A:414:ASN:ND2	3:A:620:HOH:O	2.48	0.46
1:D:394:SER:O	1:D:398:LEU:HG	2.17	0.45
1:F:397:SER:HA	1:F:410:GLN:NE2	2.31	0.45
1:F:229:PRO:HA	1:F:232:TRP:CE2	2.52	0.45
1:B:366:THR:O	1:B:370:ARG:HB3	2.15	0.45
1:A:196:VAL:HG12	1:A:200:LYS:HE3	1.99	0.45
1:E:51:ALA:HA	1:E:55:ASP:O	2.17	0.45
1:D:229:PRO:HA	1:D:232:TRP:CE2	2.52	0.45
1:F:68:GLY:HA3	1:F:332:ASN:O	2.17	0.45
1:F:15:LYS:HE3	1:F:111:GLU:OE2	2.17	0.44
1:B:229:PRO:HA	1:B:232:TRP:CE2	2.52	0.44
1:B:370:ARG:O	3:B:603:HOH:O	2.21	0.44
1:C:68:GLY:HA3	1:C:332:ASN:O	2.17	0.44
1:A:79:ILE:HD12	1:A:81:PRO:HD3	1.99	0.44
1:A:53:THR:O	1:B:288:GLU:HG2	2.18	0.44
1:F:406:THR:O	1:F:409:MET:HG2	2.17	0.44
1:F:152:GLN:NE2	3:F:607:HOH:O	2.35	0.44
1:F:379:LEU:HD22	1:F:431:CYS:HB2	2.00	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:258:PHE:HD2	1:B:340:TRP:CH2	2.35	0.44
1:E:122:LEU:HD21	1:E:126:PRO:HD3	1.98	0.44
1:B:292:ALA:O	1:B:295:LYS:HE3	2.18	0.44
1:A:182:GLY:HA2	3:A:652:HOH:O	2.18	0.43
1:A:46:LYS:HB2	1:A:46:LYS:HE2	1.74	0.43
1:C:78:GLU:OE1	1:C:102:LYS:NZ	2.36	0.43
1:D:40:PRO:HG2	1:D:43:LEU:HB3	2.01	0.43
1:A:175:LYS:HE2	3:A:601:HOH:O	2.18	0.43
1:B:437:GLU:O	1:B:440:LYS:HD2	2.17	0.43
1:F:375:TYR:CZ	1:F:429:VAL:HG22	2.54	0.43
1:A:455:GLU:HA	1:A:458:LYS:HB3	2.01	0.43
1:A:92:PHE:CD2	1:D:175:LYS:HD2	2.54	0.43
1:C:229:PRO:HA	1:C:232:TRP:CE2	2.54	0.43
1:E:129:TRP:CD1	1:E:248:PRO:HB2	2.54	0.43
1:E:378:ILE:HD13	1:E:459:VAL:HG21	2.00	0.43
1:C:122:LEU:HD21	1:C:126:PRO:HD3	2.00	0.42
1:D:109:ALA:HA	1:D:302:VAL:HA	2.00	0.42
1:D:89:LEU:HD12	1:D:94:TRP:CZ2	2.54	0.42
1:B:43:LEU:HA	1:B:46:LYS:HB2	2.02	0.42
1:E:2:ILE:HD13	1:E:2:ILE:HA	1.78	0.42
1:F:202:LYS:HE2	1:F:202:LYS:HB3	1.84	0.42
1:E:12:ASN:ND2	1:E:14:ASP:OD1	2.51	0.42
1:A:158:TRP:CD1	1:A:258:PHE:CD2	3.08	0.42
1:C:389:ASP:O	1:C:393:THR:HG23	2.20	0.42
1:D:192:LEU:HG	1:D:357:VAL:HG22	2.02	0.42
1:E:136:ASP:HA	1:E:146:ALA:HB2	2.02	0.41
1:C:426:PHE:O	1:C:431:CYS:HB3	2.20	0.41
1:D:79:ILE:O	1:D:79:ILE:HG13	2.20	0.41
1:C:209:ASP:HB2	3:C:739:HOH:O	2.20	0.41
1:E:349:ASN:OD1	1:E:354:ARG:NH1	2.51	0.41
1:F:384:PHE:O	1:F:387:MET:HG2	2.20	0.41
1:C:46:LYS:HB2	1:C:46:LYS:HE2	1.77	0.41
1:E:130:GLU:O	1:E:133:PRO:HD2	2.19	0.41
1:A:426:PHE:O	1:A:431:CYS:HB3	2.20	0.41
1:E:152:GLN:HA	1:E:348:ILE:HD11	2.02	0.41
1:F:435:LEU:HD12	1:F:438:LEU:HD11	2.02	0.41
1:F:433:ASP:OD1	1:F:456:LYS:HE2	2.21	0.41
1:E:455:GLU:O	1:E:459:VAL:HG12	2.20	0.41
1:C:109:ALA:HA	1:C:302:VAL:HA	2.02	0.41
1:E:435:LEU:HD23	1:E:452:LEU:HD11	2.02	0.41
1:C:374:GLU:O	1:C:378:ILE:HD12	2.21	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:88:LYS:O	1:C:305:LYS:HG3	2.21	0.41
1:E:179:LYS:HA	1:E:179:LYS:HD3	1.87	0.41
1:B:151:LEU:HD11	1:B:204:MET:HE3	2.03	0.41
1:A:371:ALA:HB1	1:A:374:GLU:HG3	2.03	0.40
1:E:229:PRO:HA	1:E:232:TRP:CE2	2.56	0.40
1:A:158:TRP:CE2	1:A:258:PHE:CE2	3.09	0.40
1:E:62:TRP:CD1	1:E:66:ARG:HG3	2.57	0.40
1:A:224:MET:HE3	1:A:224:MET:HB2	1.97	0.40
1:E:423:GLU:O	1:E:427:GLN:HA	2.22	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	459/477 (96%)	452 (98%)	7 (2%)	0	100	100
1	B	460/477 (96%)	453 (98%)	7 (2%)	0	100	100
1	C	462/477 (97%)	452 (98%)	10 (2%)	0	100	100
1	D	459/477 (96%)	449 (98%)	10 (2%)	0	100	100
1	E	462/477 (97%)	450 (97%)	12 (3%)	0	100	100
1	F	459/477 (96%)	446 (97%)	13 (3%)	0	100	100
All	All	2761/2862 (96%)	2702 (98%)	59 (2%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains [i](#)

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	368/384 (96%)	367 (100%)	1 (0%)	94	96
1	B	369/384 (96%)	363 (98%)	6 (2%)	70	73
1	C	371/384 (97%)	367 (99%)	4 (1%)	80	83
1	D	369/384 (96%)	364 (99%)	5 (1%)	74	77
1	E	370/384 (96%)	364 (98%)	6 (2%)	70	73
1	F	369/384 (96%)	363 (98%)	6 (2%)	70	73
All	All	2216/2304 (96%)	2188 (99%)	28 (1%)	78	79

All (28) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	258	PHE
1	B	233	SER
1	B	238	SER
1	B	258	PHE
1	B	295	LYS
1	B	440	LYS
1	B	461	LYS
1	C	258	PHE
1	C	440	LYS
1	C	451	ASN
1	C	461	LYS
1	D	25	LYS
1	D	258	PHE
1	D	309	GLU
1	D	344[A]	ARG
1	D	344[B]	ARG
1	E	34	LYS
1	E	55	ASP
1	E	258	PHE
1	E	397	SER
1	E	455	GLU
1	E	457	SER
1	F	227	ASN
1	F	258	PHE
1	F	344[A]	ARG
1	F	344[B]	ARG
1	F	427	GLN

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Mol	Chain	Res	Type
1	F	438	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	72	GLN
1	B	253	GLN
1	C	86	GLN
1	E	450	ASN
1	F	152	GLN
1	F	253	GLN

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](#)

6 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	MAL	A	501	-	24,24,24	0.50	0	35,35,35	0.57	0
2	MAL	B	501	-	24,24,24	0.49	0	35,35,35	0.57	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MAL	C	501	-	24,24,24	0.47	0	35,35,35	0.60	0
2	MAL	D	501	-	24,24,24	0.52	0	35,35,35	0.61	0
2	MAL	E	501	-	24,24,24	0.51	0	35,35,35	0.72	0
2	MAL	F	501	-	24,24,24	0.48	0	35,35,35	0.58	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	MAL	A	501	-	-	0/8/48/48	0/2/2/2
2	MAL	B	501	-	-	0/8/48/48	0/2/2/2
2	MAL	C	501	-	-	0/8/48/48	0/2/2/2
2	MAL	D	501	-	-	0/8/48/48	0/2/2/2
2	MAL	E	501	-	-	0/8/48/48	0/2/2/2
2	MAL	F	501	-	-	0/8/48/48	0/2/2/2

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
2	C	501	MAL	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	460/477 (96%)	0.30	6 (1%) 79 80	22, 31, 47, 62	0
1	B	462/477 (96%)	0.47	11 (2%) 62 63	20, 31, 60, 75	0
1	C	461/477 (96%)	0.43	14 (3%) 54 55	24, 33, 54, 74	0
1	D	459/477 (96%)	0.44	13 (2%) 56 57	23, 35, 56, 84	0
1	E	463/477 (97%)	0.59	25 (5%) 29 31	23, 38, 56, 77	0
1	F	459/477 (96%)	0.87	43 (9%) 11 11	26, 42, 76, 90	0
All	All	2764/2862 (96%)	0.52	112 (4%) 41 42	20, 35, 59, 90	0

All (112) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	459	VAL	5.6
1	F	239	ALA	5.5
1	F	453	ARG	5.3
1	F	443	PRO	4.7
1	F	375	TYR	4.4
1	F	450	ASN	4.4
1	E	137	LYS	4.3
1	B	460	ALA	4.3
1	E	401	TYR	4.2
1	F	449	VAL	4.1
1	E	460	ALA	4.0
1	E	120	ASP	3.9
1	F	457	SER	3.8
1	F	441	ASP	3.8
1	D	451	ASN	3.8
1	F	123	PRO	3.7
1	E	459	VAL	3.7
1	F	447	ASN	3.6
1	E	124	ASN	3.6

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Mol	Chain	Res	Type	RSRZ
1	B	459	VAL	3.5
1	E	122	LEU	3.5
1	F	390	TYR	3.4
1	C	451	ASN	3.3
1	F	124	ASN	3.3
1	B	413	TYR	3.3
1	D	258	PHE	3.3
1	F	235	ILE	3.3
1	C	445	LEU	3.3
1	F	391	HIS	3.2
1	B	458	LYS	3.2
1	D	354	ARG	3.2
1	E	457	SER	3.2
1	E	139	LEU	3.2
1	A	454	LYS	3.1
1	E	461	LYS	3.1
1	C	447	ASN	3.1
1	F	120	ASP	3.1
1	F	458	LYS	3.0
1	E	121	LEU	3.0
1	F	141	ALA	2.9
1	F	341	TYR	2.8
1	F	440	LYS	2.8
1	F	218	ASN	2.8
1	A	459	VAL	2.8
1	E	231	ALA	2.8
1	E	123	PRO	2.8
1	E	134	ALA	2.7
1	E	241	ASN	2.7
1	A	458	LYS	2.7
1	C	448	LEU	2.7
1	D	375	TYR	2.6
1	D	455	GLU	2.6
1	F	223	ALA	2.6
1	B	258	PHE	2.6
1	C	457	SER	2.6
1	E	463	ILE	2.6
1	D	450	ASN	2.6
1	B	457	SER	2.5
1	F	413	TYR	2.5
1	A	258	PHE	2.5
1	D	458	LYS	2.5

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Mol	Chain	Res	Type	RSRZ
1	A	447	ASN	2.5
1	F	435	LEU	2.5
1	F	452	LEU	2.5
1	F	431	CYS	2.5
1	C	460	ALA	2.5
1	D	459	VAL	2.4
1	E	441	ASP	2.4
1	F	6	LYS	2.4
1	E	2	ILE	2.4
1	D	453	ARG	2.4
1	D	454	LYS	2.4
1	B	96	ALA	2.4
1	C	52	ALA	2.4
1	C	441	ASP	2.3
1	E	258	PHE	2.3
1	C	455	GLU	2.3
1	C	390	TYR	2.3
1	F	1	LYS	2.3
1	C	2	ILE	2.3
1	E	235	ILE	2.3
1	C	450	ASN	2.3
1	D	341	TYR	2.2
1	E	295	LYS	2.2
1	E	144	LYS	2.2
1	C	33	ILE	2.2
1	F	222	THR	2.2
1	F	438	LEU	2.2
1	F	456	LYS	2.2
1	F	117	TYR	2.2
1	F	269	ALA	2.2
1	F	119	LYS	2.2
1	F	238	SER	2.2
1	F	79	ILE	2.1
1	D	1	LYS	2.1
1	F	17	TYR	2.1
1	F	34	LYS	2.1
1	F	433	ASP	2.1
1	F	122	LEU	2.1
1	B	97	VAL	2.1
1	F	125	PRO	2.1
1	E	119	LYS	2.1
1	B	123	PRO	2.1

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Mol	Chain	Res	Type	RSRZ
1	A	390	TYR	2.1
1	E	50	VAL	2.0
1	F	371	ALA	2.0
1	F	232	TRP	2.0
1	F	454	LYS	2.0
1	D	378	ILE	2.0
1	B	53	THR	2.0
1	B	341	TYR	2.0
1	E	135	LEU	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
2	MAL	A	501	23/23	0.93	0.13	-0.30	23,29,33,35	0
2	MAL	F	501	23/23	0.89	0.14	-0.37	28,35,41,46	0
2	MAL	D	501	23/23	0.93	0.13	-0.93	23,30,35,38	0
2	MAL	C	501	23/23	0.94	0.12	-0.94	23,30,33,36	0
2	MAL	E	501	23/23	0.90	0.13	-0.94	28,33,35,39	0
2	MAL	B	501	23/23	0.95	0.12	-1.34	20,26,30,32	0

6.5 Other polymers [i](#)

There are no such residues in this entry.