



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 19, 2025 – 04:00 PM EDT

PDB ID : 4GAO
Title : DCNL complex with N-terminally acetylated NEDD8 E2 peptide
Authors : Monda, J.K.; Scott, D.C.; Miller, D.J.; Harper, J.W.; Bennett, E.J.; Schulman, B.A.
Deposited on : 2012-07-25
Resolution : 3.28 Å(reported)

This is a Full wwPDB X-ray Structure Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.21
EDS	:	3.0
Percentile statistics	:	20231227.v01 (using entries in the PDB archive December 27th 2023)
CCP4	:	9.0.004 (Gargrove)
Density-Fitness	:	1.0.11
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.41.4

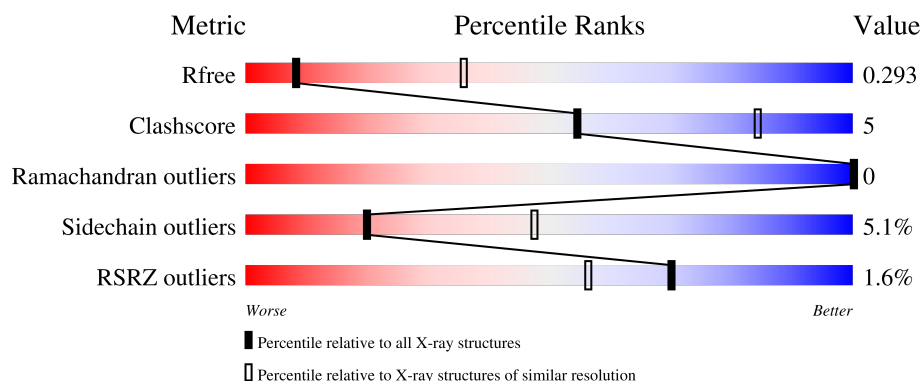
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 3.28 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}	164625	1214 (3.30-3.26)
Clashscore	180529	1265 (3.30-3.26)
Ramachandran outliers	177936	1264 (3.30-3.26)
Sidechain outliers	177891	1263 (3.30-3.26)
RSRZ outliers	164620	1215 (3.30-3.26)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	200	<div> <div>2%</div> <div> <div></div> <div>80%</div> <div>16%</div> <div>.</div> </div> </div>
1	B	200	<div> <div>%</div> <div> <div></div> <div>84%</div> <div>11%</div> <div>.</div> </div> </div>
1	D	200	<div> <div>2%</div> <div> <div></div> <div>81%</div> <div>12%</div> <div>.</div> <div>6%</div> </div> </div>
1	G	200	<div> <div>%</div> <div> <div></div> <div>82%</div> <div>14%</div> <div>.</div> </div> </div>
2	C	12	<div> <div></div> <div> <div>75%</div> <div>8%</div> <div>17%</div> </div> </div>

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Mol	Chain	Length	Quality of chain
2	E	12	<div><div></div><div>8%</div><div>75%</div><div>8%</div><div>17%</div></div>
2	F	12	<div><div></div><div>83%</div><div>8%</div><div>8%</div></div>
2	H	12	<div><div></div><div>75%</div><div>8%</div><div>17%</div></div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 6299 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 DCN1-like protein 2.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	193	Total	C	N	O	S	0	0	0
			1511	978	240	284	9			
1	B	192	Total	C	N	O	S	0	0	0
			1501	972	236	284	9			
1	D	188	Total	C	N	O	S	0	0	0
			1504	974	238	283	9			
1	G	193	Total	C	N	O	S	0	0	0
			1511	974	241	287	9			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	60	GLY	-	expression tag	UNP Q6PH85
A	61	SER	-	expression tag	UNP Q6PH85
B	60	GLY	-	expression tag	UNP Q6PH85
B	61	SER	-	expression tag	UNP Q6PH85
D	60	GLY	-	expression tag	UNP Q6PH85
D	61	SER	-	expression tag	UNP Q6PH85
G	60	GLY	-	expression tag	UNP Q6PH85
G	61	SER	-	expression tag	UNP Q6PH85

- Molecule 2 is a protein called NEDD8-conjugating enzyme Ubc12.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	10	Total	C	N	O	S	0	0	0
			66	43	11	11	1			
2	E	10	Total	C	N	O	S	0	0	0
			69	46	10	12	1			
2	F	11	Total	C	N	O	S	0	0	0
			67	43	11	12	1			
2	H	10	Total	C	N	O	S	0	0	0
			68	46	10	11	1			

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	1	AME	MET	acetylation	UNP P61081
E	1	AME	MET	acetylation	UNP P61081
F	1	AME	MET	acetylation	UNP P61081
H	1	AME	MET	acetylation	UNP P61081

- Molecule 3 is BROMIDE ION (three-letter code: BR) (formula: Br).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total Br 1 1	0	0
3	B	1	Total Br 1 1	0	0

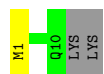
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- Molecule 1: DCN1-like protein 2




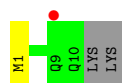
- Molecule 2: NEDD8-conjugating enzyme Ubc12

Chain C:  75% 8% 17%




- Molecule 2: NEDD8-conjugating enzyme Ubc12

Chain E:  8% 75% 8% 17%



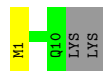
- Molecule 2: NEDD8-conjugating enzyme Ubc12

Chain F:  83% 8% 8%



- Molecule 2: NEDD8-conjugating enzyme Ubc12

Chain H:  75% 8% 17%



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	48.57Å 190.15Å 49.07Å 90.00° 101.75° 90.00°	Depositor
Resolution (Å)	40.00 – 3.28 40.00 – 3.29	Depositor EDS
% Data completeness (in resolution range)	98.1 (40.00-3.28) 98.1 (40.00-3.29)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.20 (at 3.25Å)	Xtriage
Refinement program	REFMAC 5.5.0102	Depositor
R, R_{free}	0.252 , 0.296 0.246 , 0.293	Depositor DCC
R_{free} test set	651 reflections (4.94%)	wwPDB-VP
Wilson B-factor (Å ²)	55.3	Xtriage
Anisotropy	0.095	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.32 , 33.0	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtriage
Estimated twinning fraction	0.070 for l,-k,h	Xtriage
F_o, F_c correlation	0.87	EDS
Total number of atoms	6299	wwPDB-VP
Average B, all atoms (Å ²)	51.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.85% 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: BR, AME

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	0.47	0/1549	0.52	0/2102
1	B	0.45	0/1538	0.49	0/2088
1	D	0.48	0/1542	0.52	0/2088
1	G	0.50	0/1548	0.51	0/2099
2	C	0.45	0/54	0.42	0/72
2	E	0.52	0/58	0.41	0/78
2	F	0.40	0/55	0.39	0/75
2	H	0.50	0/57	0.41	0/77
All	All	0.47	0/6401	0.51	0/8679

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	1511	0	1407	17	0
1	B	1501	0	1401	15	0
1	D	1504	0	1422	12	0
1	G	1511	0	1406	14	0
2	C	66	0	59	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	E	69	0	58	0	0
2	F	67	0	50	0	0
2	H	68	0	55	0	0
3	A	1	0	0	1	0
3	B	1	0	0	0	0
All	All	6299	0	5858	57	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 5.

All (57) 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:B:175:LEU:HD22	1:B:205:LYS:HA	1.50	0.91
1:G:175:LEU:HD22	1:G:205:LYS:HA	1.55	0.89
1:D:212:THR:HG21	1:D:239:LEU:HD23	1.69	0.75
1:D:141:LEU:HB2	1:D:142:PRO:HD3	1.71	0.71
1:B:183:LYS:O	1:B:187:SER:HB2	1.94	0.68
1:A:141:LEU:HB2	1:A:142:PRO:HD3	1.76	0.68
1:B:175:LEU:CD2	1:B:205:LYS:HA	2.24	0.67
1:D:162:PHE:CE2	1:D:210:ARG:HG2	2.29	0.67
1:D:140:LEU:O	1:D:144:LEU:HG	1.95	0.67
1:A:183:LYS:O	1:A:187:SER:HB2	1.98	0.63
1:G:179:VAL:O	1:G:183:LYS:HG3	1.98	0.63
1:A:179:VAL:O	1:A:183:LYS:HG3	2.03	0.58
1:D:144:LEU:O	1:D:147:GLU:HG2	2.03	0.57
1:A:141:LEU:CB	1:A:142:PRO:HD3	2.38	0.54
1:D:93:LEU:HG	1:D:134:MET:CE	2.37	0.54
1:A:93:LEU:HG	1:A:134:MET:CE	2.38	0.54
1:A:182:TRP:CD1	1:A:186:LEU:HD12	2.43	0.54
1:G:98:ALA:O	1:G:184:LEU:HD21	2.10	0.52
1:A:95:LEU:HD11	1:A:141:LEU:HD11	1.91	0.52
1:D:105:ILE:HD11	1:D:141:LEU:HD21	1.93	0.50
1:A:105:ILE:HD11	1:A:141:LEU:HD21	1.95	0.47
1:G:175:LEU:CD2	1:G:205:LYS:HA	2.38	0.46
1:G:196:TRP:HA	1:G:243:PHE:CE1	2.50	0.46
1:G:114:GLN:HA	1:G:164:PHE:CZ	2.51	0.46
1:G:122:PHE:HD2	1:G:123:LEU:HD12	1.80	0.46
1:B:141:LEU:HB2	1:B:142:PRO:HD3	1.98	0.46
1:A:108:LYS:HD2	1:A:108:LYS:HA	1.82	0.45
1:G:141:LEU:HB2	1:G:142:PRO:HD3	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:183:LYS:O	1:G:187:SER:HB2	2.16	0.45
1:A:237:PRO:HD2	1:A:240:ILE:HD12	1.99	0.45
1:B:223:ILE:HA	1:B:229:ASN:HD21	1.81	0.45
1:A:230:TYR:CE1	1:A:236:TRP:CD1	3.05	0.44
1:G:94:SER:HB2	1:G:138:LYS:NZ	2.33	0.44
1:G:167:ASN:O	1:G:170:GLN:HB2	2.17	0.44
1:G:204:HIS:O	1:G:204:HIS:ND1	2.50	0.44
1:A:120:LYS:HB2	3:A:301:BR:BR	2.72	0.44
1:B:140:LEU:O	1:B:144:LEU:HG	2.18	0.44
1:B:167:ASN:N	1:B:167:ASN:OD1	2.50	0.44
1:D:93:LEU:HG	1:D:134:MET:HE1	1.99	0.43
1:A:124:ASP:O	1:A:128:GLU:HG3	2.18	0.43
1:A:136:LYS:O	1:A:140:LEU:HD13	2.19	0.43
1:D:191:LYS:HD2	1:D:192:PHE:CE2	2.53	0.43
1:B:175:LEU:HB2	1:B:206:ARG:O	2.19	0.43
1:B:167:ASN:HD21	1:D:150:ASP:CG	2.22	0.42
1:B:248:ARG:HB2	1:B:249:PRO:HD3	2.01	0.42
1:B:120:LYS:HB3	1:B:120:LYS:HE3	1.86	0.42
1:B:230:TYR:HB3	1:B:248:ARG:HH21	1.84	0.42
1:D:114:GLN:HA	1:D:164:PHE:CZ	2.55	0.42
1:A:64:LYS:HB3	1:A:132:ASP:HB2	2.01	0.41
1:A:196:TRP:HA	1:A:243:PHE:CE1	2.56	0.41
1:D:100:ILE:O	1:D:104:VAL:HG23	2.19	0.41
1:G:202:GLU:HB3	1:G:203:HIS:CD2	2.56	0.41
1:A:197:ASN:O	1:A:201:MET:HG3	2.20	0.41
1:B:232:GLU:C	1:B:234:GLY:H	2.24	0.41
1:B:179:VAL:O	1:B:183:LYS:HG3	2.21	0.40
1:B:222:MET:O	1:B:229:ASN:ND2	2.55	0.40
1:G:135:GLU:H	1:G:135:GLU:HG3	1.52	0.40

There are no symmetry-related clashes.

5.3 Torsion angles ⓘ

5.3.1 Protein backbone ⓘ

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	191/200 (96%)	182 (95%)	9 (5%)	0	100	100
1	B	190/200 (95%)	184 (97%)	6 (3%)	0	100	100
1	D	186/200 (93%)	182 (98%)	4 (2%)	0	100	100
1	G	191/200 (96%)	185 (97%)	6 (3%)	0	100	100
2	C	8/12 (67%)	8 (100%)	0	0	100	100
2	E	8/12 (67%)	8 (100%)	0	0	100	100
2	F	9/12 (75%)	9 (100%)	0	0	100	100
2	H	8/12 (67%)	8 (100%)	0	0	100	100
All	All	791/848 (93%)	766 (97%)	25 (3%)	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	153/175 (87%)	146 (95%)	7 (5%)	23	51
1	B	153/175 (87%)	147 (96%)	6 (4%)	27	55
1	D	156/175 (89%)	146 (94%)	10 (6%)	14	40
1	G	154/175 (88%)	145 (94%)	9 (6%)	17	43
2	C	3/11 (27%)	3 (100%)	0	100	100
2	E	4/11 (36%)	4 (100%)	0	100	100
2	F	2/11 (18%)	2 (100%)	0	100	100
2	H	3/11 (27%)	3 (100%)	0	100	100
All	All	628/744 (84%)	596 (95%)	32 (5%)	20	47

All (32) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	66	GLU
1	A	110	ARG

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Mol	Chain	Res	Type
1	A	187	SER
1	A	191	LYS
1	A	210	ARG
1	A	217	LEU
1	A	222	MET
1	B	145	GLU
1	B	150	ASP
1	B	167	ASN
1	B	215	LEU
1	B	217	LEU
1	B	222	MET
1	D	119	ARG
1	D	135	GLU
1	D	150	ASP
1	D	163	THR
1	D	187	SER
1	D	191	LYS
1	D	210	ARG
1	D	215	LEU
1	D	217	LEU
1	D	222	MET
1	G	73	LYS
1	G	134	MET
1	G	135	GLU
1	G	149	LYS
1	G	160	PHE
1	G	163	THR
1	G	187	SER
1	G	217	LEU
1	G	222	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	203	HIS
1	G	159	GLN
1	G	203	HIS

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues 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	AME	C	1	2	9,10,11	0.41	0	9,11,13	1.04	1 (11%)
2	AME	H	1	2	9,10,11	0.60	0	9,11,13	1.09	1 (11%)
2	AME	F	1	2	9,10,11	0.40	0	9,11,13	1.35	1 (11%)
2	AME	E	1	2	9,10,11	0.50	0	9,11,13	1.13	1 (11%)

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	AME	C	1	2	-	0/9/10/12	-
2	AME	H	1	2	-	0/9/10/12	-
2	AME	F	1	2	-	0/9/10/12	-
2	AME	E	1	2	-	2/9/10/12	-

There are no bond length outliers.

All (4) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	F	1	AME	CB-CA-N	-3.09	104.89	110.52
2	H	1	AME	CB-CA-N	-2.56	105.85	110.52
2	C	1	AME	CB-CA-N	-2.18	106.55	110.52
2	E	1	AME	CB-CA-N	-2.08	106.73	110.52

There are no chirality outliers.

All (2) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	E	1	AME	CA-CB-CG-SD
2	E	1	AME	C-CA-CB-CG

There are no ring outliers.

No monomer is involved in short contacts.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 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 [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	193/200 (96%)	0.23	5 (2%) 57 42	25, 50, 83, 90	0
1	B	192/200 (96%)	0.21	2 (1%) 79 67	25, 48, 76, 87	0
1	D	188/200 (94%)	0.26	3 (1%) 70 56	27, 55, 84, 96	0
1	G	193/200 (96%)	0.16	2 (1%) 79 67	28, 44, 80, 92	0
2	C	9/12 (75%)	0.68	0 100 100	44, 47, 50, 50	0
2	E	9/12 (75%)	0.37	1 (11%) 12 11	38, 46, 55, 56	0
2	F	10/12 (83%)	0.06	0 100 100	40, 48, 55, 55	0
2	H	9/12 (75%)	0.65	0 100 100	39, 46, 52, 52	0
All	All	803/848 (94%)	0.22	13 (1%) 70 56	25, 49, 81, 96	0

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	60	GLY	3.5
1	A	233	GLU	3.4
1	B	251	VAL	2.8
1	B	188	GLY	2.6
1	D	62	LYS	2.5
1	G	60	GLY	2.4
1	A	79	ASN	2.4
1	D	208	ILE	2.3
2	E	9	GLN	2.3
1	A	232	GLU	2.3
1	D	235	ALA	2.2
1	G	252	THR	2.1
1	A	171	LYS	2.0

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

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
2	AME	C	1	11/12	0.93	0.11	43,43,44,44	0
2	AME	H	1	11/12	0.94	0.10	39,39,40,40	0
2	AME	F	1	11/12	0.95	0.09	38,38,39,39	0
2	AME	E	1	11/12	0.95	0.10	35,36,37,37	0

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å ²)	Q<0.9
3	BR	B	301	1/1	0.98	0.04	52,52,52,52	0
3	BR	A	301	1/1	0.99	0.08	46,46,46,46	0

6.5 Other polymers [i](#)

There are no such residues in this entry.