



Full wwPDB X-ray Structure Validation Report ⓘ

May 13, 2020 – 08:16 am BST

PDB ID : 1U2O
Title : Crystal Structure Of The N-Domain Of Grp94 Lacking The Charged Domain
In Complex With Neca
Authors : Soldano, K.L.; Jivan, A.; Nicchitta, C.V.; Gewirth, D.T.
Deposited on : 2004-07-19
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

<https://www.wwpdb.org/validation/2017/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 : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.11
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.11

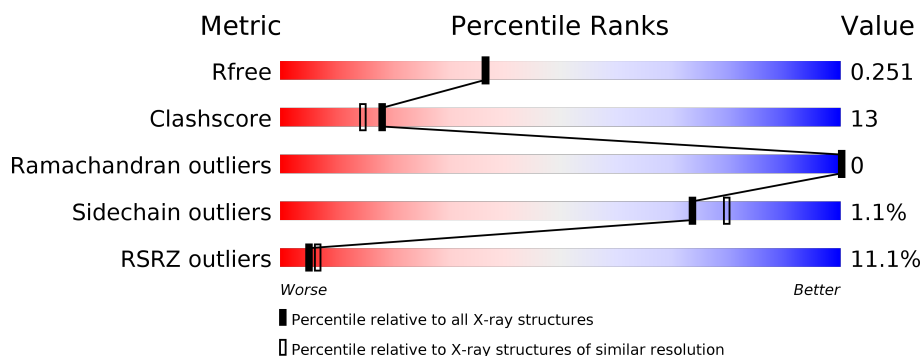
1 Overall quality at a glance

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}	130704	5197 (2.10-2.10)
Clashscore	141614	5710 (2.10-2.10)
Ramachandran outliers	138981	5647 (2.10-2.10)
Sidechain outliers	138945	5648 (2.10-2.10)
RSRZ outliers	127900	5083 (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 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	236	<div> <div>11%</div> <div>69%</div> <div>27%</div> <div>.</div> </div>
1	B	236	<div> <div>11%</div> <div>78%</div> <div>18%</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	PG4	A	2002	-	-	X	-
3	PG4	B	2009	-	-	X	-
4	1PE	B	2006	-	-	X	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 4003 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.

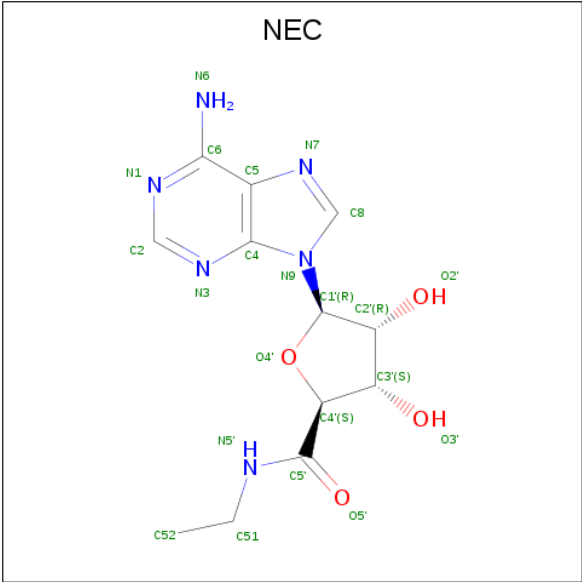
- Molecule 1 is a protein called Endoplasmin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	226	Total	C	N	O	S	0	3	0
			1768	1116	300	345	7			
1	B	226	Total	C	N	O	S	0	7	0
			1783	1125	299	351	8			

There are 8 discrepancies between the modelled and reference sequences:

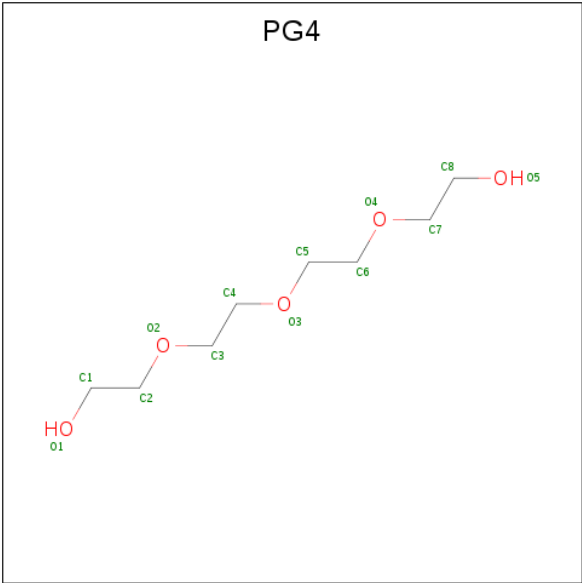
Chain	Residue	Modelled	Actual	Comment	Reference
A	65	GLY	-	CLONING ARTIFACT	UNP P41148
A	66	SER	-	CLONING ARTIFACT	UNP P41148
A	67	HIS	-	CLONING ARTIFACT	UNP P41148
A	68	MET	-	CLONING ARTIFACT	UNP P41148
B	65	GLY	-	CLONING ARTIFACT	UNP P41148
B	66	SER	-	CLONING ARTIFACT	UNP P41148
B	67	HIS	-	CLONING ARTIFACT	UNP P41148
B	68	MET	-	CLONING ARTIFACT	UNP P41148

- Molecule 2 is N-ETHYL-5'-CARBOXAMIDO ADENOSINE (three-letter code: NEC) (formula: C₁₂H₁₆N₆O₄).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	A	1	Total	C	N	O	0	0
			22	12	6	4		
2	B	1	Total	C	N	O	0	0
			22	12	6	4		

- Molecule 3 is TETRAETHYLENE GLYCOL (three-letter code: PG4) (formula: C₈H₁₈O₅).



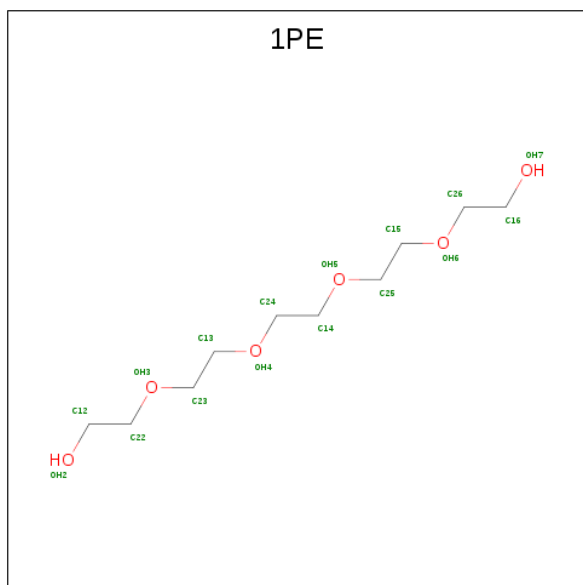
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			10	6	4		
3	A	1	Total	C	O	0	0
			13	8	5		

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	B	1	Total	C	O	0	0
			13	8	5		
3	B	1	Total	C	O	0	0
			13	8	5		
3	B	1	Total	C	O	0	0
			7	4	3		

- Molecule 4 is PENTAETHYLENE GLYCOL (three-letter code: 1PE) (formula: $C_{10}H_{22}O_6$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	B	1	Total	C	O	0	0
			16	10	6		

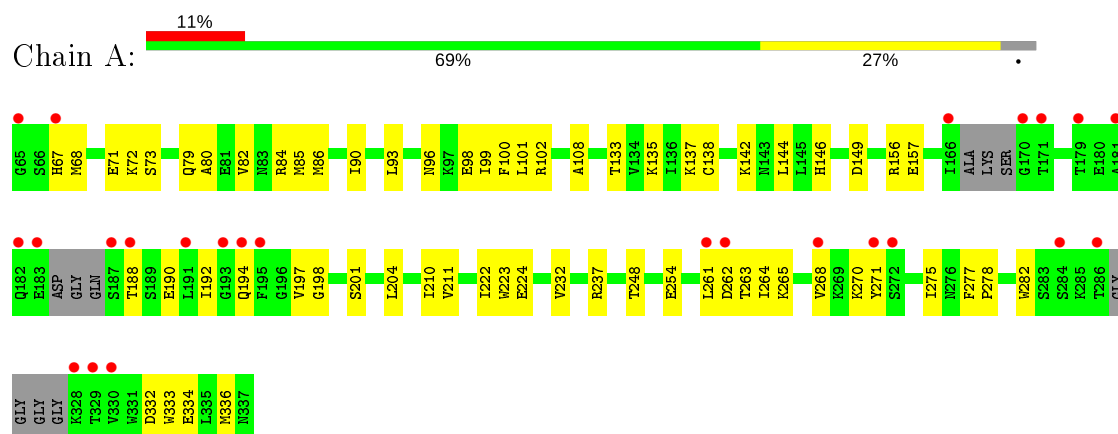
- Molecule 5 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	175	Total	O	0	0
			175	175		
5	B	161	Total	O	0	0
			161	161		

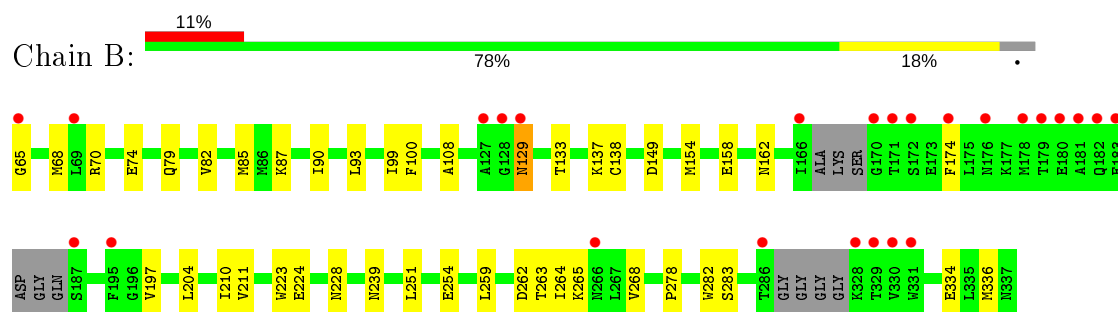
3 Residue-property plots [i](#)

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



• Molecule 1: Endoplasmin



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	65.30Å 84.20Å 94.09Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 2.10 47.05 – 2.10	Depositor EDS
% Data completeness (in resolution range)	87.1 (50.00-2.10) 86.9 (47.05-2.10)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.05	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.03 (at 2.10Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.213 , 0.257 0.208 , 0.251	Depositor DCC
R_{free} test set	2683 reflections (9.83%)	wwPDB-VP
Wilson B-factor (Å ²)	29.6	Xtriage
Anisotropy	0.168	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.33 , 64.4	EDS
L-test for twinning ²	$\langle L \rangle = 0.45$, $\langle L^2 \rangle = 0.28$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4003	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 34.18 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.1041e-04. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: PG4, NEC, 1PE

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	0.33	0/1791	0.58	0/2417
1	B	0.33	0/1808	0.59	0/2445
All	All	0.33	0/3599	0.59	0/4862

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

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	1768	0	1719	46	0
1	B	1783	0	1695	42	0
2	A	22	0	16	0	0
2	B	22	0	16	1	0
3	A	23	0	31	12	0
3	B	33	0	45	13	0
4	B	16	0	22	9	0
5	A	175	0	0	2	0
5	B	161	0	0	3	0
All	All	4003	0	3544	97	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 13.

All (97) 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:137:LYS:HZ2	3:B:2009:PG4:H42	1.12	1.12
1:B:87:LYS:HE3	4:B:2006:1PE:H251	1.46	0.95
3:A:2001:PG4:O1	3:A:2001:PG4:H31	1.83	0.77
1:A:137:LYS:HZ3	3:A:2002:PG4:H41	1.48	0.77
1:B:137:LYS:NZ	3:B:2009:PG4:H42	1.99	0.74
1:B:85[A]:MET:SD	1:B:197:VAL:HG11	2.28	0.74
1:B:282:TRP:CZ2	3:B:2009:PG4:H82	2.29	0.67
1:A:86:MET:HB3	4:B:2006:1PE:H231	1.76	0.67
1:B:137:LYS:NZ	3:B:2009:PG4:H21	2.13	0.63
1:A:137:LYS:HE3	5:A:2086:HOH:O	2.02	0.59
1:A:68:MET:HG3	1:A:73:SER:OG	2.02	0.59
3:B:2008:PG4:H32	3:B:2009:PG4:H11	1.85	0.58
1:B:282:TRP:CH2	3:B:2009:PG4:H82	2.38	0.58
1:A:137:LYS:HZ3	3:A:2002:PG4:H51	1.67	0.58
1:B:133:THR:HG22	1:B:278:PRO:HG2	1.84	0.58
1:B:87:LYS:CE	4:B:2006:1PE:H251	2.30	0.56
1:B:65:GLY:HA3	5:B:2142:HOH:O	2.06	0.56
1:A:67:HIS:CG	1:A:68:MET:N	2.74	0.56
1:A:90:ILE:HA	1:A:204:LEU:HD21	1.88	0.55
1:B:70:ARG:HD2	5:B:2123:HOH:O	2.06	0.55
1:B:254[B]:GLU:N	1:B:254[B]:GLU:OE1	2.37	0.54
1:B:137:LYS:NZ	3:B:2009:PG4:H52	2.23	0.54
1:B:137:LYS:HZ2	3:B:2009:PG4:H21	1.72	0.53
1:A:85:MET:SD	1:A:197:VAL:HG11	2.48	0.53
3:B:2008:PG4:C3	3:B:2009:PG4:H11	2.38	0.53
1:A:237:ARG:HH22	3:A:2001:PG4:H42	1.73	0.53
1:B:211:VAL:HB	1:B:223:TRP:HB3	1.91	0.52
1:A:275:ILE:HG22	1:A:277:PHE:H	1.74	0.52
3:A:2001:PG4:O1	3:A:2001:PG4:C3	2.56	0.51
1:B:87:LYS:NZ	4:B:2006:1PE:OH2	2.42	0.51
1:A:282:TRP:CH2	3:A:2002:PG4:H11	2.47	0.50
1:B:334:GLU:O	1:B:336:MET:HE2	2.10	0.50
1:A:102:ARG:HB2	1:A:271:TYR:CD1	2.47	0.50
1:B:265:LYS:HE2	1:B:283:SER:OG	2.10	0.50
1:B:99:ILE:HD12	1:B:99:ILE:C	2.31	0.50
1:B:251:LEU:HD12	1:B:259:LEU:HG	1.93	0.50
1:A:137:LYS:NZ	3:A:2002:PG4:H22	2.28	0.49
1:A:262[B]:ASP:OD1	1:A:263:THR:N	2.46	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:138:CYS:SG	1:A:264:ILE:HG13	2.52	0.49
1:A:190:GLU:O	1:A:194:GLN:HG3	2.12	0.48
1:A:264:ILE:O	1:A:268:VAL:HG23	2.13	0.48
1:B:210:ILE:HD12	1:B:210:ILE:N	2.27	0.48
3:A:2001:PG4:HO1	3:A:2001:PG4:H31	1.75	0.48
1:B:90:ILE:HD12	4:B:2006:1PE:H161	1.94	0.48
1:A:210:ILE:HB	1:A:248:THR:HB	1.95	0.48
1:B:138:CYS:SG	1:B:264:ILE:HG13	2.54	0.48
1:A:282:TRP:CZ2	3:A:2002:PG4:H11	2.49	0.48
1:B:262[A]:ASP:OD2	1:B:263:THR:N	2.48	0.47
1:A:86:MET:CB	4:B:2006:1PE:H231	2.42	0.46
1:B:87:LYS:HE2	4:B:2006:1PE:H221	1.97	0.46
1:B:158:GLU:O	1:B:162:ASN:HB2	2.16	0.45
3:B:2009:PG4:H42	3:B:2009:PG4:H21	1.68	0.45
1:B:129:ASN:OD1	3:B:2011:PG4:H41	2.16	0.45
1:A:84:ARG:HG2	1:B:228:ASN:ND2	2.32	0.45
1:B:108:ALA:HB1	1:B:149:ASP:HB3	1.99	0.45
3:B:2009:PG4:H52	3:B:2009:PG4:H71	1.67	0.45
1:A:72:LYS:HA	5:A:2091:HOH:O	2.15	0.45
1:A:192:ILE:CG2	1:A:198:GLY:HA2	2.48	0.44
3:A:2002:PG4:H22	3:A:2002:PG4:H41	1.68	0.44
1:A:99:ILE:C	1:A:99:ILE:HD12	2.38	0.44
1:A:334:GLU:O	1:A:336:MET:HE2	2.17	0.44
1:B:90:ILE:HD12	4:B:2006:1PE:C16	2.48	0.44
1:A:108:ALA:HB1	1:A:149:ASP:HB3	2.00	0.44
1:A:237:ARG:HH12	3:A:2001:PG4:C4	2.31	0.44
1:A:261:LEU:O	1:A:265:LYS:HG3	2.18	0.43
1:A:93:LEU:HD23	1:A:99:ILE:HD11	2.00	0.43
1:A:156[A]:ARG:HG2	1:A:157:GLU:OE2	2.19	0.43
1:A:254:GLU:N	1:A:254:GLU:OE1	2.42	0.43
1:A:96:ASN:OD1	1:A:98[A]:GLU:OE1	2.36	0.43
1:B:85[B]:MET:HE1	1:B:174:PHE:HD2	1.82	0.43
1:B:68:MET:HG2	1:B:74:GLU:OE1	2.18	0.42
1:A:211:VAL:HB	1:A:223:TRP:HB3	2.01	0.42
1:B:137:LYS:HZ1	3:B:2009:PG4:H21	1.83	0.42
1:A:222:ILE:O	1:A:232:VAL:HA	2.20	0.42
1:B:264:ILE:O	1:B:268:VAL:HG23	2.19	0.42
1:A:80:ALA:O	1:A:84:ARG:HG3	2.20	0.42
1:A:67:HIS:O	1:A:71:GLU:HG2	2.20	0.42
1:B:79:GLN:HB3	1:B:82:VAL:HG23	2.00	0.42
1:B:85[B]:MET:CE	1:B:174:PHE:HD2	2.33	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:90:ILE:HA	1:B:204:LEU:CD2	2.50	0.42
1:A:79:GLN:HB3	1:A:82:VAL:HG23	2.01	0.41
1:B:99:ILE:HD12	1:B:100:PHE:N	2.35	0.41
1:A:100:PHE:CG	1:A:101:LEU:N	2.88	0.41
1:B:154:MET:SD	2:B:1002:NEC:H1'	2.60	0.41
1:A:100:PHE:HA	1:A:201:SER:OG	2.21	0.41
1:A:282:TRP:CZ3	1:A:332:ASP:HA	2.56	0.41
1:B:90:ILE:HA	1:B:204:LEU:HD21	2.02	0.41
1:A:270:LYS:O	1:A:270:LYS:HG2	2.20	0.41
1:A:135:LYS:HB3	1:A:333:TRP:CZ3	2.57	0.40
1:A:142:LYS:HB3	1:A:144:LEU:HG	2.03	0.40
1:B:251:LEU:HD12	1:B:259:LEU:CD2	2.52	0.40
1:B:239:ASN:HB2	5:B:2075:HOH:O	2.21	0.40
1:A:137:LYS:HG2	1:A:146:HIS:HB2	2.03	0.40
1:A:188:THR:O	1:A:192:ILE:HG12	2.21	0.40
1:A:133:THR:HG22	1:A:278:PRO:HG2	2.04	0.40
4:B:2006:1PE:H122	4:B:2006:1PE:H232	1.79	0.40
3:A:2002:PG4:H72	3:A:2002:PG4:H51	1.66	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	221/236 (94%)	210 (95%)	11 (5%)	0	100	100
1	B	225/236 (95%)	217 (96%)	8 (4%)	0	100	100
All	All	446/472 (94%)	427 (96%)	19 (4%)	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	189/207 (91%)	188 (100%)	1 (0%)	88	92
1	B	189/207 (91%)	186 (98%)	3 (2%)	62	69
All	All	378/414 (91%)	374 (99%)	4 (1%)	73	79

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	224	GLU
1	B	93	LEU
1	B	129	ASN
1	B	224	GLU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

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
3	PG4	B	2009	-	12,12,12	0.52	0	11,11,11	1.29	1 (9%)
3	PG4	B	2008	-	12,12,12	0.50	0	11,11,11	0.63	0
4	1PE	B	2006	-	15,15,15	0.53	0	14,14,14	1.18	0
3	PG4	A	2002	-	12,12,12	0.52	0	11,11,11	1.34	1 (9%)
3	PG4	B	2011	-	6,6,12	0.50	0	5,5,11	0.37	0
3	PG4	A	2001	-	9,9,12	0.53	0	8,8,11	0.35	0
2	NEC	A	1001	-	21,24,24	2.57	6 (28%)	21,35,35	1.57	2 (9%)
2	NEC	B	1002	-	21,24,24	2.55	5 (23%)	21,35,35	1.46	2 (9%)

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	PG4	B	2009	-	-	4/10/10/10	-
3	PG4	B	2008	-	-	0/10/10/10	-
4	1PE	B	2006	-	-	4/13/13/13	-
3	PG4	A	2002	-	-	4/10/10/10	-
3	PG4	B	2011	-	-	1/4/4/10	-
3	PG4	A	2001	-	-	4/7/7/10	-
2	NEC	A	1001	-	-	3/7/27/27	0/3/3/3
2	NEC	B	1002	-	-	3/7/27/27	0/3/3/3

All (11) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	1002	NEC	O5'-C5'	9.21	1.41	1.23
2	A	1001	NEC	O5'-C5'	9.09	1.41	1.23
2	A	1001	NEC	C5'-N5'	4.98	1.44	1.33
2	B	1002	NEC	C5'-N5'	4.75	1.44	1.33
2	A	1001	NEC	C4-N3	2.75	1.39	1.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1001	NEC	O4'-C4'	-2.57	1.40	1.43
2	B	1002	NEC	C4-N3	2.47	1.39	1.35
2	B	1002	NEC	O4'-C4'	-2.26	1.40	1.43
2	B	1002	NEC	C2-N3	2.16	1.35	1.32
2	A	1001	NEC	C2-N3	2.12	1.35	1.32
2	A	1001	NEC	O4'-C1'	2.09	1.44	1.41

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	1001	NEC	O4'-C1'-C2'	-4.44	100.44	106.93
2	B	1002	NEC	O4'-C1'-C2'	-4.32	100.62	106.93
2	A	1001	NEC	C1'-O4'-C4'	3.64	115.42	109.11
2	B	1002	NEC	C1'-O4'-C4'	3.15	114.58	109.11
3	A	2002	PG4	C5-O3-C4	-3.09	99.88	113.29
3	B	2009	PG4	C5-O3-C4	-3.02	100.21	113.29

There are no chirality outliers.

All (23) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1001	NEC	O4'-C4'-C5'-N5'
2	B	1002	NEC	O4'-C4'-C5'-N5'
3	A	2001	PG4	C3-C4-O3-C5
4	B	2006	1PE	C24-C14-OH5-C25
2	A	1001	NEC	O4'-C4'-C5'-O5'
2	B	1002	NEC	O4'-C4'-C5'-O5'
3	B	2009	PG4	C4-C3-O2-C2
3	A	2002	PG4	C5-C6-O4-C7
2	A	1001	NEC	C3'-C4'-C5'-N5'
2	B	1002	NEC	C3'-C4'-C5'-N5'
3	B	2009	PG4	C5-C6-O4-C7
4	B	2006	1PE	C25-C15-OH6-C26
3	B	2011	PG4	C1-C2-O2-C3
3	A	2002	PG4	C4-C3-O2-C2
4	B	2006	1PE	C12-C22-OH3-C23
3	A	2001	PG4	C4-C3-O2-C2
3	A	2001	PG4	C1-C2-O2-C3
3	A	2002	PG4	O3-C5-C6-O4
3	B	2009	PG4	O2-C3-C4-O3
3	B	2009	PG4	O3-C5-C6-O4
3	A	2002	PG4	O2-C3-C4-O3

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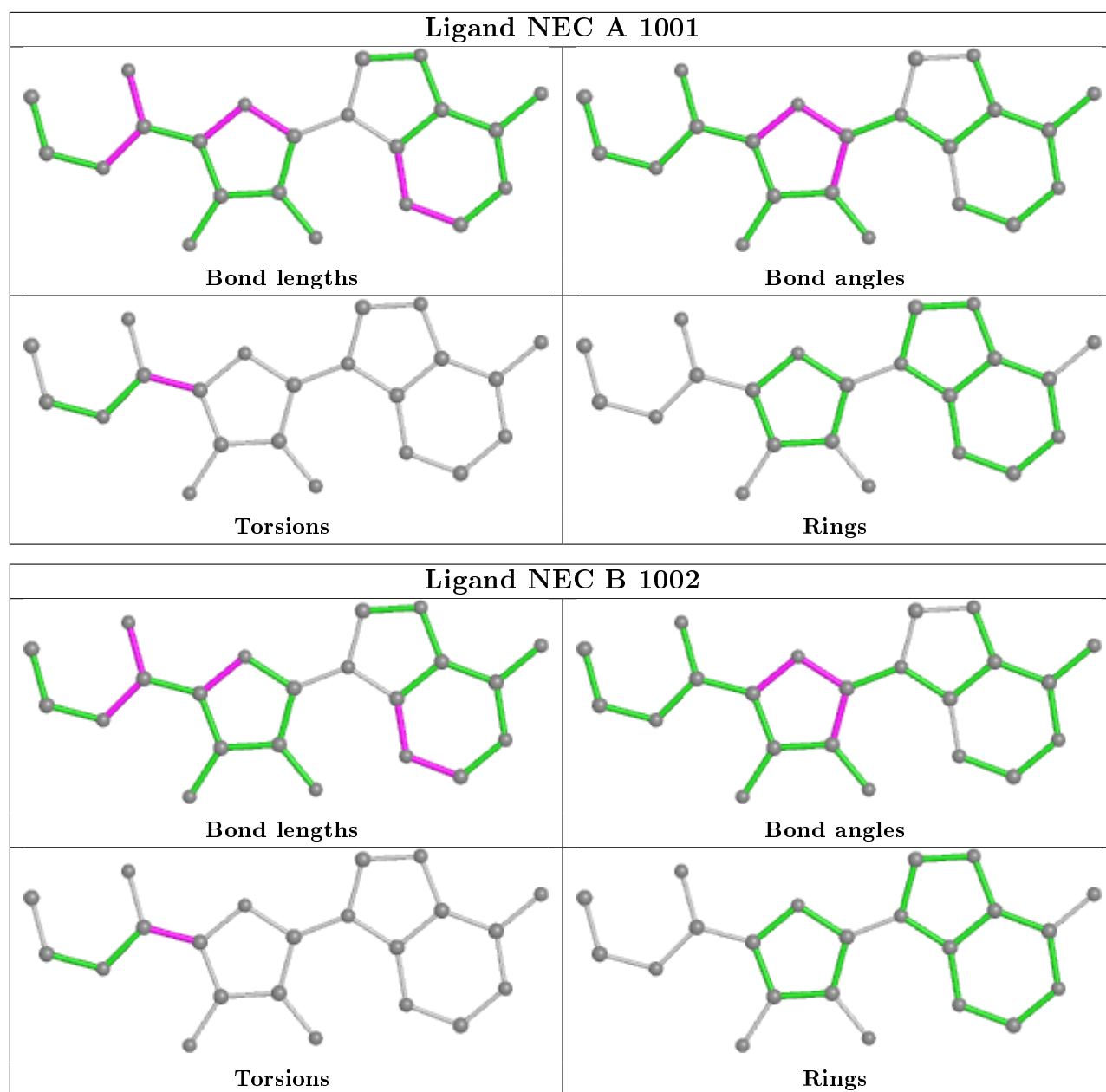
Mol	Chain	Res	Type	Atoms
4	B	2006	1PE	OH6-C15-C25-OH5
3	A	2001	PG4	O2-C3-C4-O3

There are no ring outliers.

7 monomers are involved in 35 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	B	2009	PG4	12	0
3	B	2008	PG4	2	0
4	B	2006	1PE	9	0
3	A	2002	PG4	7	0
3	B	2011	PG4	1	0
3	A	2001	PG4	5	0
2	B	1002	NEC	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



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	226/236 (95%)	0.56	25 (11%) 5 7	17, 32, 74, 89	0
1	B	226/236 (95%)	0.61	25 (11%) 5 7	17, 29, 77, 91	0
All	All	452/472 (95%)	0.59	50 (11%) 5 7	17, 30, 76, 91	0

All (50) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	181	ALA	8.1
1	B	330	VAL	6.0
1	B	174	PHE	5.8
1	B	187	SER	5.8
1	B	172	SER	5.6
1	B	182	GLN	5.5
1	B	179	THR	5.4
1	B	170	GLY	5.3
1	A	182	GLN	5.3
1	B	286	THR	4.9
1	A	181	ALA	4.7
1	A	286	THR	4.4
1	B	183	GLU	4.4
1	A	179	THR	4.3
1	A	271	TYR	4.0
1	B	195	PHE	3.8
1	B	129	ASN	3.8
1	A	195	PHE	3.6
1	A	171	THR	3.6
1	A	170	GLY	3.5
1	A	330	VAL	3.5
1	B	128	GLY	3.5
1	A	261	LEU	3.4
1	B	176	ASN	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	65	GLY	3.2
1	A	272	SER	3.1
1	A	191	LEU	3.0
1	B	65	GLY	2.8
1	A	187	SER	2.8
1	B	180	GLU	2.8
1	B	69	LEU	2.7
1	A	329	THR	2.7
1	B	166	ILE	2.7
1	A	183	GLU	2.7
1	A	328	LYS	2.6
1	A	268	VAL	2.6
1	B	127	ALA	2.5
1	A	67	HIS	2.5
1	A	284	SER	2.4
1	A	166	ILE	2.4
1	A	194	GLN	2.3
1	A	193	GLY	2.3
1	A	262[A]	ASP	2.2
1	B	329	THR	2.2
1	A	188	THR	2.2
1	B	328	LYS	2.2
1	B	266[A]	ASN	2.2
1	B	178	MET	2.1
1	B	171	THR	2.1
1	B	331	TRP	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](#)

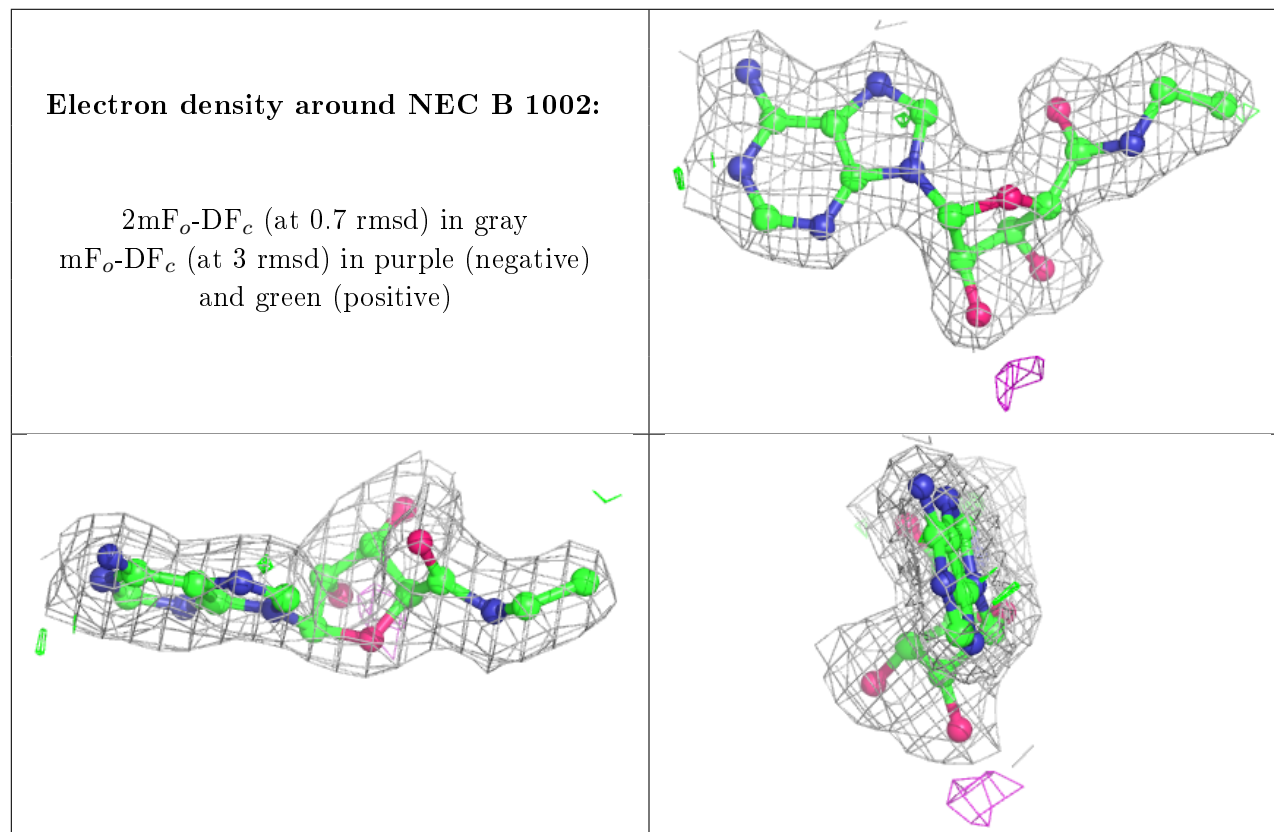
There are no carbohydrates in this entry.

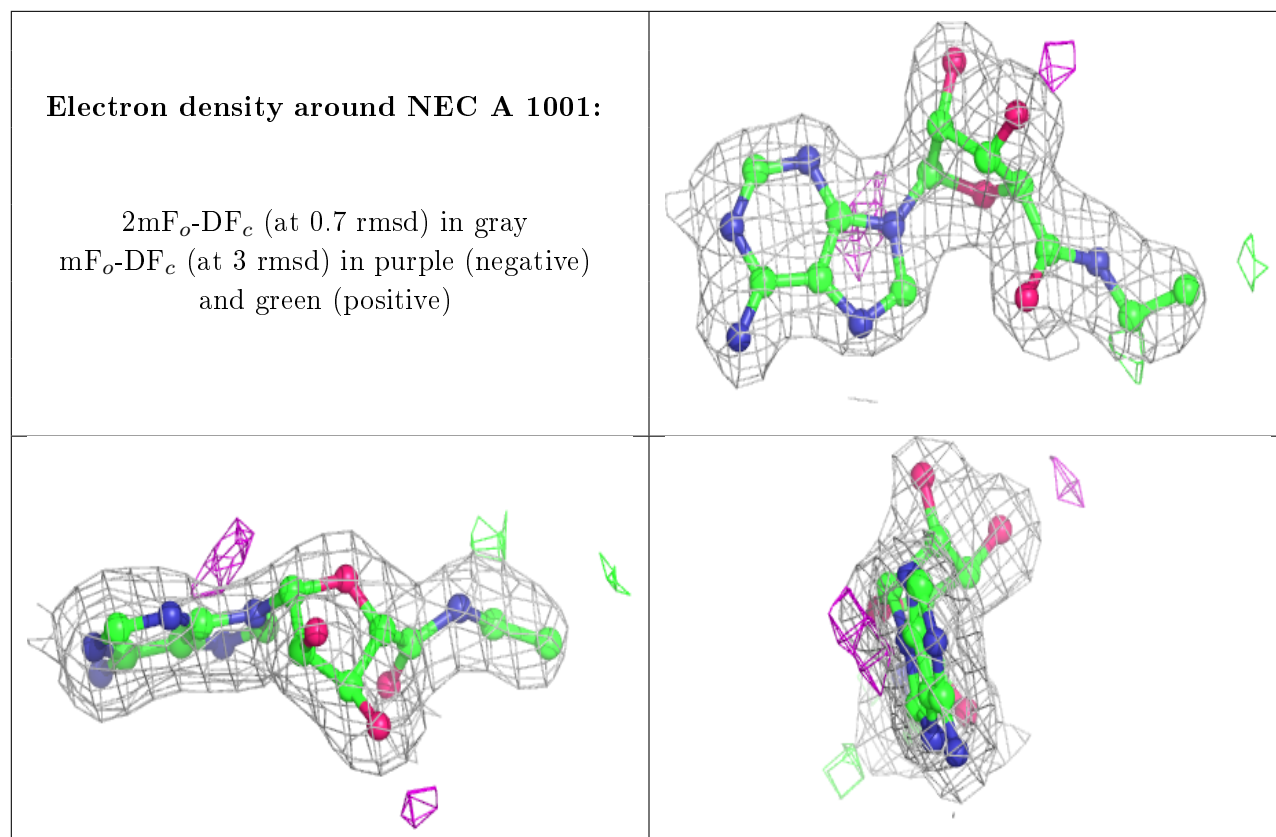
6.4 Ligands [i](#)

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. 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(\AA^2)	Q<0.9
3	PG4	B	2009	13/13	0.66	0.24	51,55,59,59	0
3	PG4	A	2001	10/13	0.73	0.28	37,41,44,49	0
3	PG4	A	2002	13/13	0.75	0.25	59,61,63,64	0
4	1PE	B	2006	16/16	0.80	0.35	46,49,58,60	0
3	PG4	B	2011	7/13	0.87	0.22	61,61,62,63	0
3	PG4	B	2008	13/13	0.88	0.20	39,43,50,52	0
2	NEC	B	1002	22/22	0.91	0.15	20,28,32,33	0
2	NEC	A	1001	22/22	0.93	0.15	17,27,35,38	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





6.5 Other polymers [i](#)

There are no such residues in this entry.