Full wwPDB X-ray Structure Validation Report (i)

Jun 22, 2024 – 10:22 AM EDT

PDB ID	:	6IAG
Title	:	Crystal structure of human phosphodiesterase 4D2 catalytic domain with inhibitor NPD-637
Authors	:	Singh, A.K.; Brown, D.G.
Deposited on	:	2018-11-26
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
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) 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 (i)) were used in the production of this report:

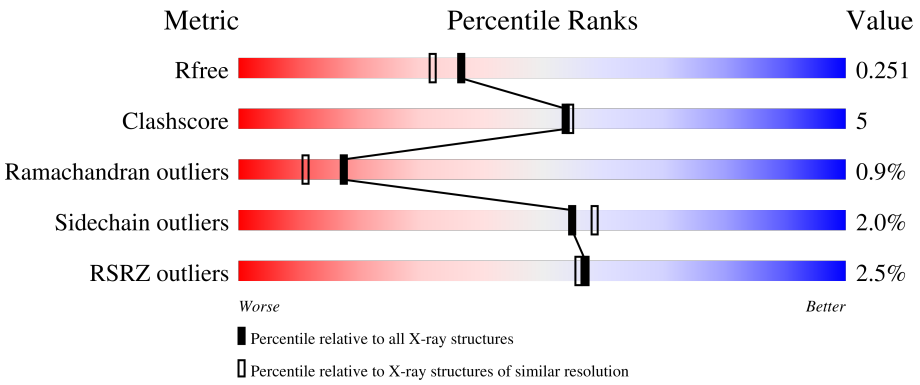
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

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}	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (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 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 <=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	364	5%, 81%, 8%, • , 10%
1	B	364	81%, 7%, • , 11%
1	C	364	3%, 77%, 10%, • , 11%
1	D	364	%, 78%, 10%, • , 11%

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
5	E3Q	A	513	-	-	-	X

2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 11530 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 cAMP-specific 3',5'-cyclic phosphodiesterase 4D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	329	2658	1681	454	509	14	0	0	0
1	B	324	2622	1659	448	501	14	0	0	0
1	C	323	2613	1654	446	499	14	0	0	0
1	D	324	2628	1663	449	502	14	0	1	0

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	75	GLY	-	expression tag	UNP Q08499
A	76	SER	-	expression tag	UNP Q08499
A	77	HIS	-	expression tag	UNP Q08499
A	78	MET	-	expression tag	UNP Q08499
B	75	GLY	-	expression tag	UNP Q08499
B	76	SER	-	expression tag	UNP Q08499
B	77	HIS	-	expression tag	UNP Q08499
B	78	MET	-	expression tag	UNP Q08499
C	75	GLY	-	expression tag	UNP Q08499
C	76	SER	-	expression tag	UNP Q08499
C	77	HIS	-	expression tag	UNP Q08499
C	78	MET	-	expression tag	UNP Q08499
D	75	GLY	-	expression tag	UNP Q08499
D	76	SER	-	expression tag	UNP Q08499
D	77	HIS	-	expression tag	UNP Q08499
D	78	MET	-	expression tag	UNP Q08499

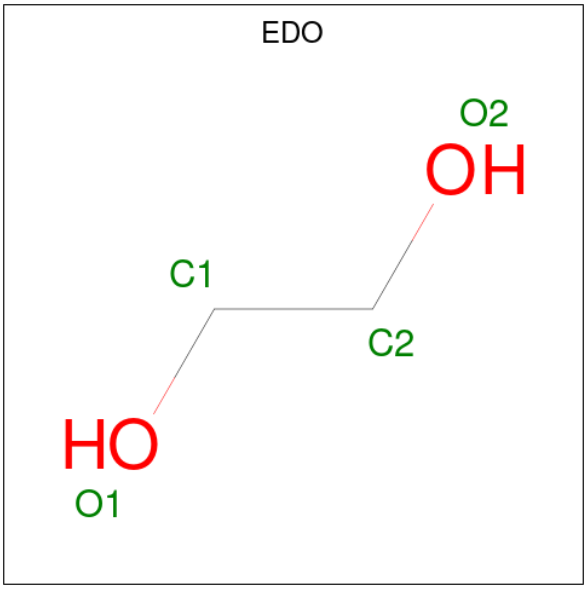
- Molecule 2 is ZINC ION (three-letter code: ZN) (formula: Zn).

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

- Molecule 3 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
4	A	1	4	2	2	0	0
			Total	C	O		
4	A	1	4	2	2	0	0

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O			
4	A	1	4	2	2		0	0
			Total	C	O			
4	A	1	4	2	2		0	0
			Total	C	O			
4	A	1	4	2	2		0	0
			Total	C	O			
4	A	1	4	2	2		0	0
			Total	C	O			
4	A	1	4	2	2		0	0
			Total	C	O			
4	A	1	4	2	2		0	0
			Total	C	O			
4	A	1	4	2	2		0	0
			Total	C	O			
4	A	1	4	2	2		0	0
			Total	C	O			
4	B	1	4	2	2		0	0
			Total	C	O			
4	B	1	4	2	2		0	0
			Total	C	O			
4	B	1	4	2	2		0	0
			Total	C	O			
4	B	1	4	2	2		0	0
			Total	C	O			
4	B	1	4	2	2		0	0
			Total	C	O			
4	B	1	4	2	2		0	0
			Total	C	O			
4	C	1	4	2	2		0	0
			Total	C	O			
4	C	1	4	2	2		0	0
			Total	C	O			
4	C	1	4	2	2		0	0
			Total	C	O			
4	C	1	4	2	2		0	0
			Total	C	O			
4	C	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0

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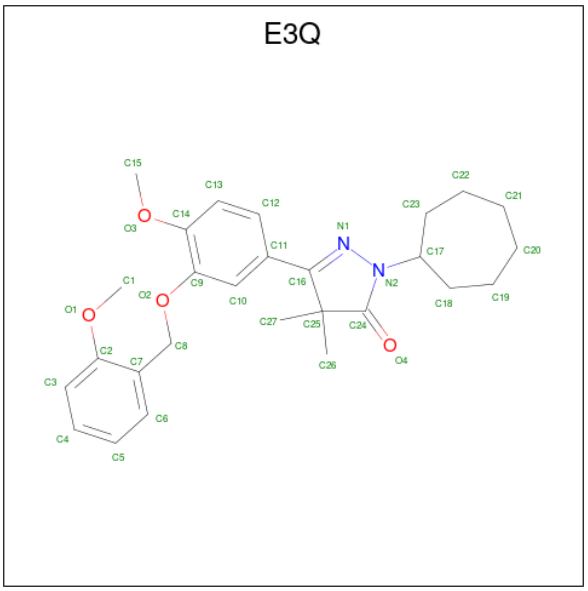
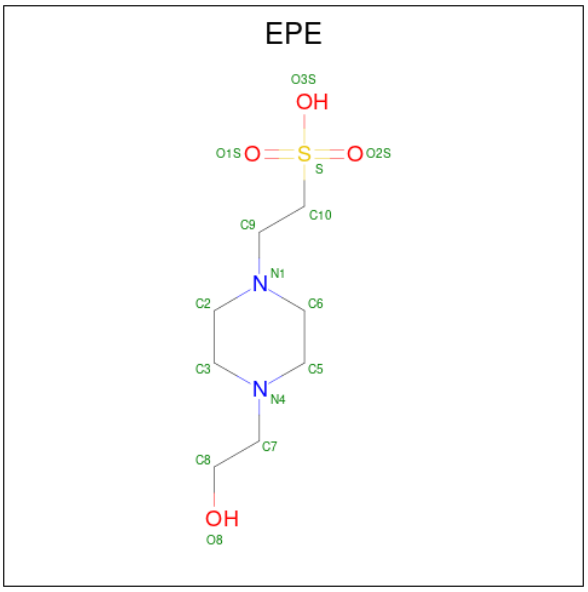
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0
			Total	C	O			
4	D	1	4	2	2		0	0

- Molecule 5 is 1-cycloheptyl-3-{4-methoxy-3-[(2-methoxyphenyl)methoxy]phenyl}-4,4-dimethyl-4,5-dihydro- 1H-pyrazol-5-one (three-letter code: E3Q) (formula: $C_{27}H_{34}N_2O_4$) (labeled as "Ligand of Interest" by depositor).



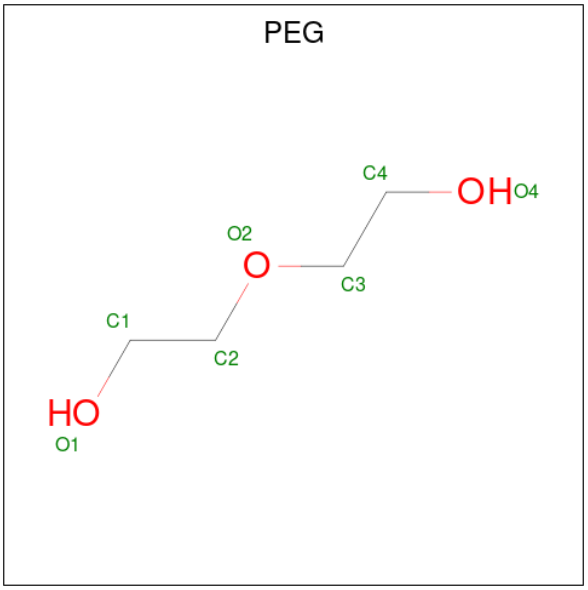
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
5	A	1	33	27	2	4	0	0
5	B	1	33	27	2	4	0	0
5	C	1	33	27	2	4	0	0
5	D	1	33	27	2	4	0	0

- Molecule 6 is 4-(2-HYDROXYETHYL)-1-PIPERAZINE ETHANESULFONIC ACID (three-letter code: EPE) (formula: $C_8H_{18}N_2O_4S$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	N	O	S		
6	B	1	15	8	2	4	1	0	0
			Total	C	N	O	S		
6	B	1	15	8	2	4	1	0	0
			Total	C	N	O	S		
6	C	1	15	8	2	4	1	0	0
			Total	C	N	O	S		
6	D	1	15	8	2	4	1	0	0

- Molecule 7 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
7	D	1	7	4	3	0	0
			Total	C	O		
7	D	1	7	4	3	0	0

- Molecule 8 is water.

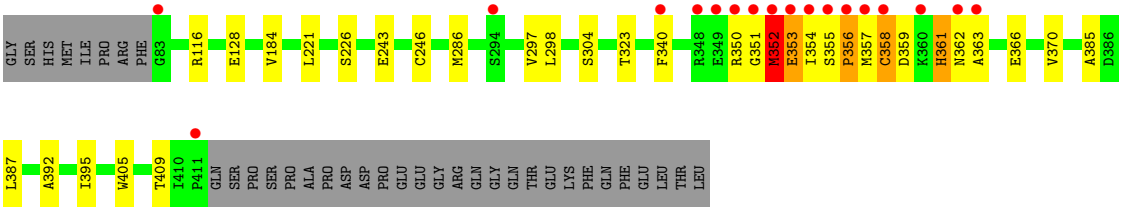
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
8	A	165	165	165	0	0
			Total	O		
8	B	161	161	161	0	0
			Total	O		
8	C	129	129	129	0	0
			Total	O		
8	D	200	200	200	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide 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.

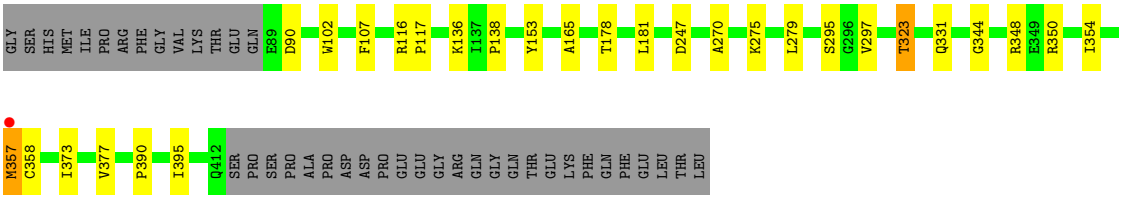
• Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D

Chain A:



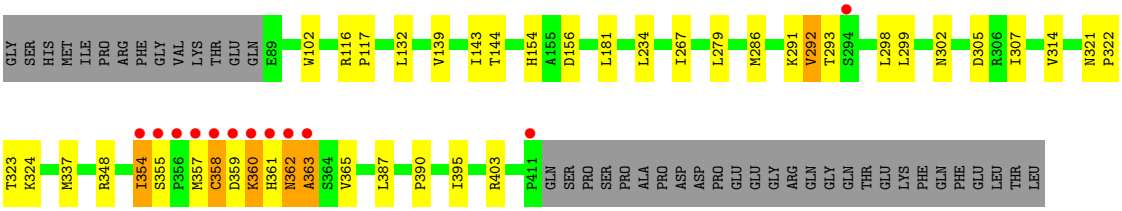
• Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D

Chain B:



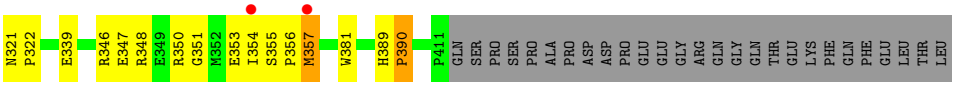
• Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D

Chain C:



• Molecule 1: cAMP-specific 3',5'-cyclic phosphodiesterase 4D

Chain D:


N321
P322
E339
R346
E347
R348
E349
R350
G351
M352
E353
I354
S355
P356
M357
W381
H389
P390
P411
GLN
SER
PRO
SER
PRO
ALA
PRO
ASP
ASP
PRO
GLU
GLU
GLY
ARG
GLN
GLY
GLN
THR
GLU
LYS
PHE
GLN
PHE
GLU
LEU
THR
LEU

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α, β, γ	99.11Å 110.31Å 160.72Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	84.50 – 2.00 84.36 – 2.00	Depositor EDS
% Data completeness (in resolution range)	98.1 (84.50-2.00) 98.1 (84.36-2.00)	Depositor EDS
R_{merge}	0.12	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.15 (at 2.00Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.200 , 0.246 0.210 , 0.251	Depositor DCC
R_{free} test set	5902 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	27.9	Xtriage
Anisotropy	0.165	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 52.8	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.95	EDS
Total number of atoms	11530	wwPDB-VP
Average B, all atoms (Å^2)	38.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.92% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ 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: E3Q, EDO, PEG, MG, ZN, EPE

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	0.77	1/2712 (0.0%)	0.80	0/3684
1	B	0.74	0/2676	0.81	0/3636
1	C	0.74	0/2667	0.80	0/3624
1	D	0.74	0/2685	0.79	0/3648
All	All	0.75	1/10740 (0.0%)	0.80	0/14592

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	116	ARG	C-N	8.48	1.50	1.34

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	2658	0	2616	24	0
1	B	2622	0	2578	22	0
1	C	2613	0	2570	34	0
1	D	2628	0	2586	32	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	1	0	0	0	0
2	D	1	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
3	C	1	0	0	0	0
3	D	1	0	0	0	0
4	A	40	0	60	1	0
4	B	24	0	36	3	0
4	C	20	0	30	0	0
4	D	56	0	84	9	0
5	A	33	0	0	2	0
5	B	33	0	0	1	0
5	C	33	0	0	0	0
5	D	33	0	0	1	0
6	B	30	0	34	3	0
6	C	15	0	18	0	0
6	D	15	0	18	1	0
7	D	14	0	20	3	0
8	A	165	0	0	0	0
8	B	161	0	0	3	0
8	C	129	0	0	0	0
8	D	200	0	0	3	0
All	All	11530	0	10650	113	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 (113) 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:A:351:GLY:O	1:A:353:GLU:N	1.61	1.31
1:A:359:ASP:O	1:A:363:ALA:HB3	1.27	1.30
1:C:359:ASP:O	1:C:361:HIS:N	1.93	1.01
1:C:358:CYS:O	1:C:359:ASP:OD1	1.79	0.99
1:B:354:ILE:HG12	8:B:602:HOH:O	1.66	0.94
1:A:359:ASP:O	1:A:363:ALA:CB	2.16	0.91
1:A:351:GLY:C	1:A:353:GLU:H	1.75	0.91
1:C:354:ILE:HD13	1:C:359:ASP:OD2	1.81	0.80
1:B:107:PHE:HB2	6:B:508:EPE:H101	1.64	0.77
1:A:361:HIS:O	1:A:363:ALA:N	2.18	0.76
1:C:354:ILE:CG2	1:C:354:ILE:O	2.35	0.74

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:346:ARG:NH2	8:D:601:HOH:O	2.21	0.74
1:D:115:ASN:HD21	4:D:505:EDO:H12	1.53	0.73
1:D:115:ASN:HD21	4:D:505:EDO:C1	2.02	0.72
1:D:174:VAL:HG12	7:D:511:PEG:H41	1.71	0.72
1:C:357:MET:O	1:C:363:ALA:HB1	1.90	0.71
1:D:286:MET:CE	1:D:308[A]:GLN:OE1	2.41	0.68
1:B:102:TRP:HE1	1:B:323:THR:HG23	1.59	0.68
1:C:358:CYS:O	1:C:359:ASP:CG	2.32	0.66
1:C:348:ARG:HH11	1:C:354:ILE:HD11	1.59	0.66
1:C:354:ILE:O	1:C:354:ILE:HG22	1.97	0.64
1:C:348:ARG:NH1	1:C:354:ILE:HD11	2.12	0.64
1:A:352:MET:O	1:A:353:GLU:HB3	1.98	0.63
1:C:354:ILE:CD1	1:C:359:ASP:OD2	2.47	0.61
1:A:243:GLU:HB2	1:A:246:CYS:SG	2.41	0.61
1:D:107:PHE:HB2	6:D:516:EPE:H91	1.83	0.60
1:C:359:ASP:C	1:C:361:HIS:H	2.04	0.60
1:D:115:ASN:ND2	4:D:505:EDO:H12	2.16	0.59
1:C:234:LEU:HB2	4:D:515:EDO:H11	1.85	0.59
1:A:298:LEU:HD21	1:A:387:LEU:HD11	1.85	0.58
1:B:348:ARG:NH2	8:B:602:HOH:O	2.36	0.57
1:B:354:ILE:N	1:B:354:ILE:HD12	2.20	0.56
1:C:362:ASN:O	1:C:363:ALA:CB	2.55	0.54
1:A:354:ILE:C	1:A:354:ILE:HD12	2.28	0.54
1:A:350:ARG:HG2	1:C:144:THR:HG23	1.90	0.53
1:C:139:VAL:O	1:C:143:ILE:HG12	2.08	0.53
1:B:116:ARG:N	1:B:117:PRO:CD	2.73	0.52
1:B:323:THR:HG23	1:B:395:ILE:HG23	1.92	0.52
1:A:323:THR:HB	1:A:395:ILE:HG23	1.92	0.52
1:D:179:PRO:HG2	7:D:512:PEG:H21	1.90	0.52
1:C:234:LEU:CB	4:D:515:EDO:H11	2.39	0.52
1:C:321:ASN:HB2	1:C:322:PRO:HD3	1.92	0.52
1:A:352:MET:O	1:A:353:GLU:CB	2.58	0.51
1:B:136:LYS:O	1:B:138:PRO:HD3	2.10	0.51
1:D:224:ASN:OD1	4:D:515:EDO:H12	2.10	0.51
1:C:362:ASN:O	1:C:363:ALA:HB3	2.11	0.51
1:B:102:TRP:HE1	1:B:323:THR:CG2	2.22	0.50
7:D:512:PEG:H21	7:D:512:PEG:O4	2.11	0.50
1:D:286:MET:HE1	1:D:308[A]:GLN:OE1	2.10	0.50
1:A:366:GLU:O	1:A:370:VAL:HG23	2.12	0.50
1:B:357:MET:HA	8:B:637:HOH:O	2.12	0.50
1:D:346:ARG:HG3	8:D:775:HOH:O	2.11	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:355:SER:O	1:A:356:PRO:C	2.51	0.49
1:C:359:ASP:C	1:C:361:HIS:N	2.56	0.49
4:B:510:EDO:H21	1:D:243:GLU:OE1	2.13	0.49
1:A:405:TRP:O	1:A:409:THR:HG23	2.13	0.48
1:B:331:GLN:HE22	6:B:508:EPE:H102	1.79	0.48
1:C:359:ASP:O	1:C:362:ASN:N	2.37	0.47
1:D:142:LEU:O	1:D:146:LEU:HG	2.14	0.47
1:A:340:PHE:HD2	5:A:513:E3Q:C5	2.27	0.47
1:D:286:MET:HE3	1:D:308[A]:GLN:OE1	2.12	0.47
1:D:321:ASN:HB2	1:D:322:PRO:HD3	1.95	0.47
1:A:351:GLY:O	1:A:353:GLU:CA	2.57	0.47
1:D:357:MET:HB2	5:D:517:E3Q:C4	2.45	0.47
1:A:355:SER:O	1:A:358:CYS:HB2	2.16	0.46
1:B:357:MET:HB3	5:B:509:E3Q:C4	2.45	0.46
1:C:354:ILE:O	1:C:354:ILE:HG23	2.14	0.46
1:B:270:ALA:HB1	1:B:279:LEU:HD11	1.97	0.46
1:C:116:ARG:N	1:C:117:PRO:CD	2.79	0.46
1:D:303:TYR:CE1	4:D:509:EDO:H12	2.51	0.46
1:C:298:LEU:HD11	1:C:387:LEU:HG	1.97	0.46
1:D:355:SER:HB3	4:D:503:EDO:H21	1.97	0.46
4:B:510:EDO:H22	1:D:152:HIS:HE1	1.81	0.45
1:A:184:VAL:CG2	1:A:297:VAL:HG13	2.47	0.45
1:A:286:MET:HG3	4:A:512:EDO:O2	2.16	0.45
1:C:181:LEU:HD21	1:C:298:LEU:HD12	1.98	0.45
1:C:292:VAL:HG23	1:C:293:THR:N	2.32	0.45
1:A:340:PHE:HB3	1:A:358:CYS:SG	2.57	0.45
1:B:178:THR:CG2	1:B:181:LEU:HD12	2.46	0.45
1:D:348:ARG:HD3	1:D:354:ILE:HD11	1.99	0.45
1:C:286:MET:CE	1:C:305:ASP:OD1	2.65	0.45
1:D:350:ARG:NH1	8:D:611:HOH:O	2.50	0.45
1:D:104:LEU:HD11	1:D:109:ILE:CD1	2.47	0.44
1:D:304:SER:O	1:D:308[B]:GLN:HG3	2.18	0.44
1:D:347:GLU:OE1	1:D:354:ILE:HA	2.18	0.44
1:C:267:ILE:HG21	1:C:314:VAL:HG21	1.99	0.44
1:A:340:PHE:CD2	5:A:513:E3Q:C5	3.01	0.43
1:B:178:THR:HG22	1:B:181:LEU:HD12	2.00	0.43
1:C:154:HIS:HB3	1:C:156:ASP:OD1	2.18	0.43
1:D:201:ASP:O	1:D:204:HIS:HB2	2.19	0.43
1:B:153:TYR:CZ	1:B:165:ALA:HB2	2.54	0.43
6:B:508:EPE:H32	6:B:508:EPE:H81	1.79	0.42
1:C:307:ILE:HD12	1:C:307:ILE:HA	1.93	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:247:ASP:OD1	1:B:247:ASP:C	2.58	0.42
1:C:337:MET:HG3	1:C:365:VAL:HG22	2.00	0.42
1:B:344:GLY:HA3	1:B:358:CYS:O	2.19	0.42
1:B:350:ARG:NH2	4:B:510:EDO:O2	2.53	0.42
1:C:279:LEU:HD23	1:C:279:LEU:HA	1.87	0.41
1:A:385:ALA:HA	1:A:392:ALA:HB3	2.03	0.41
1:B:373:ILE:HA	1:B:377:VAL:HB	2.03	0.41
1:C:323:THR:HB	1:C:395:ILE:HG23	2.02	0.41
1:C:359:ASP:O	1:C:360:LYS:C	2.55	0.41
1:D:249:PHE:CD2	4:D:518:EDO:H11	2.55	0.41
1:D:389:HIS:HA	1:D:390:PRO:HA	1.92	0.41
1:D:206:GLY:HA2	1:D:339:GLU:OE2	2.21	0.41
1:B:295:SER:OG	1:B:297:VAL:HG23	2.21	0.41
1:D:316:CYS:HB3	1:D:381:TRP:CZ2	2.55	0.41
1:D:143:ILE:HG22	1:D:147:MET:CE	2.50	0.41
1:B:350:ARG:HH11	1:B:350:ARG:HD3	1.74	0.40
1:A:352:MET:H	1:A:352:MET:HG3	1.68	0.40
1:C:102:TRP:CE2	1:C:324:LYS:HE2	2.57	0.40
1:D:219:LEU:HD23	1:D:219:LEU:HA	1.93	0.40
1:D:104:LEU:HD22	1:D:170:GLN:HG3	2.02	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
1	A	327/364 (90%)	310 (95%)	12 (4%)	5 (2%)	10 4
1	B	322/364 (88%)	317 (98%)	5 (2%)	0	100 100
1	C	321/364 (88%)	307 (96%)	11 (3%)	3 (1%)	17 11
1	D	323/364 (89%)	315 (98%)	5 (2%)	3 (1%)	17 11
All	All	1293/1456 (89%)	1249 (97%)	33 (3%)	11 (1%)	17 11

All (11) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	352	MET
1	A	362	ASN
1	C	302	ASN
1	C	360	LYS
1	A	353	GLU
1	C	363	ALA
1	D	353	GLU
1	A	361	HIS
1	A	356	PRO
1	D	351	GLY
1	D	356	PRO

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	300/331 (91%)	293 (98%)	7 (2%)	50	53
1	B	296/331 (89%)	291 (98%)	5 (2%)	60	65
1	C	295/331 (89%)	285 (97%)	10 (3%)	37	36
1	D	297/331 (90%)	295 (99%)	2 (1%)	84	88
All	All	1188/1324 (90%)	1164 (98%)	24 (2%)	55	58

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	128	GLU
1	A	221	LEU
1	A	226	SER
1	A	304	SER
1	A	352	MET
1	A	357	MET
1	A	358	CYS
1	B	90	ASP
1	B	275	LYS

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Mol	Chain	Res	Type
1	B	323	THR
1	B	357	MET
1	B	390	PRO
1	C	132	LEU
1	C	291	LYS
1	C	292	VAL
1	C	299	LEU
1	C	354	ILE
1	C	355	SER
1	C	358	CYS
1	C	362	ASN
1	C	390	PRO
1	C	403	ARG
1	D	357	MET
1	D	390	PRO

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

Mol	Chain	Res	Type
1	A	224	ASN
1	B	331	GLN
1	D	88	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 monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 53 ligands modelled in this entry, 8 are monoatomic - leaving 45 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	$\#\|Z\| > 2$
4	EDO	A	505	-	3,3,3	0.19	0	2,2,2	0.09	0
4	EDO	D	506	-	3,3,3	0.11	0	2,2,2	0.11	0
6	EPE	B	508	-	15,15,15	2.87	2 (13%)	19,20,20	1.49	3 (15%)
4	EDO	B	507	-	3,3,3	0.13	0	2,2,2	0.03	0
4	EDO	A	507	-	3,3,3	0.15	0	2,2,2	0.29	0
4	EDO	D	518	-	3,3,3	0.30	0	2,2,2	0.52	0
4	EDO	D	507	-	3,3,3	1.07	0	2,2,2	0.80	0
6	EPE	B	511	-	15,15,15	1.85	1 (6%)	19,20,20	1.34	3 (15%)
6	EPE	D	516	-	15,15,15	1.99	1 (6%)	19,20,20	5.04	5 (26%)
4	EDO	A	509	-	3,3,3	0.55	0	2,2,2	0.36	0
4	EDO	D	508	-	3,3,3	0.23	0	2,2,2	0.42	0
4	EDO	B	506	-	3,3,3	0.25	0	2,2,2	0.53	0
6	EPE	C	505	-	15,15,15	2.43	1 (6%)	19,20,20	1.41	4 (21%)
4	EDO	D	515	-	3,3,3	0.38	0	2,2,2	0.24	0
4	EDO	A	503	-	3,3,3	0.02	0	2,2,2	0.06	0
4	EDO	C	509	-	3,3,3	0.09	0	2,2,2	0.14	0
4	EDO	D	519	-	3,3,3	0.31	0	2,2,2	0.29	0
4	EDO	C	504	-	3,3,3	0.16	0	2,2,2	0.18	0
4	EDO	A	511	-	3,3,3	0.11	0	2,2,2	0.13	0
4	EDO	D	509	-	3,3,3	0.13	0	2,2,2	0.18	0
4	EDO	C	503	-	3,3,3	0.13	0	2,2,2	0.26	0
4	EDO	B	503	-	3,3,3	0.11	0	2,2,2	0.26	0
4	EDO	D	504	-	3,3,3	0.05	0	2,2,2	0.08	0
4	EDO	A	512	-	3,3,3	0.06	0	2,2,2	0.18	0
4	EDO	B	510	-	3,3,3	0.21	0	2,2,2	0.19	0
4	EDO	D	514	-	3,3,3	0.19	0	2,2,2	0.26	0
4	EDO	D	503	-	3,3,3	0.24	0	2,2,2	0.18	0
4	EDO	D	505	-	3,3,3	0.10	0	2,2,2	0.40	0
4	EDO	D	520	-	3,3,3	0.12	0	2,2,2	0.21	0
4	EDO	D	510	-	3,3,3	0.38	0	2,2,2	0.04	0
7	PEG	D	512	-	6,6,6	0.38	0	5,5,5	0.50	0
4	EDO	C	508	-	3,3,3	0.23	0	2,2,2	0.19	0
4	EDO	B	505	-	3,3,3	0.16	0	2,2,2	0.31	0
4	EDO	D	513	-	3,3,3	0.12	0	2,2,2	0.23	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
5	E3Q	D	517	-	35,36,36	0.85	2 (5%)	43,51,51	1.41	4 (9%)
5	E3Q	C	507	-	35,36,36	0.73	2 (5%)	43,51,51	1.45	6 (13%)
4	EDO	A	506	-	3,3,3	0.20	0	2,2,2	0.41	0
4	EDO	B	504	-	3,3,3	0.19	0	2,2,2	0.13	0
7	PEG	D	511	-	6,6,6	0.28	0	5,5,5	0.37	0
5	E3Q	A	513	-	35,36,36	0.78	1 (2%)	43,51,51	1.64	6 (13%)
4	EDO	A	510	-	3,3,3	0.25	0	2,2,2	0.38	0
4	EDO	C	506	-	3,3,3	0.16	0	2,2,2	0.14	0
4	EDO	A	504	-	3,3,3	0.72	0	2,2,2	0.48	0
5	E3Q	B	509	-	35,36,36	0.88	2 (5%)	43,51,51	1.45	5 (11%)
4	EDO	A	508	-	3,3,3	0.09	0	2,2,2	0.16	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
4	EDO	A	505	-	-	1/1/1/1	-
4	EDO	D	506	-	-	0/1/1/1	-
6	EPE	B	508	-	-	6/9/19/19	0/1/1/1
4	EDO	B	507	-	-	1/1/1/1	-
4	EDO	A	507	-	-	1/1/1/1	-
4	EDO	D	518	-	-	0/1/1/1	-
4	EDO	D	507	-	-	1/1/1/1	-
6	EPE	B	511	-	-	4/9/19/19	0/1/1/1
6	EPE	D	516	-	-	1/9/19/19	0/1/1/1
4	EDO	A	509	-	-	0/1/1/1	-
4	EDO	D	508	-	-	1/1/1/1	-
4	EDO	B	506	-	-	1/1/1/1	-
6	EPE	C	505	-	-	5/9/19/19	0/1/1/1
4	EDO	D	515	-	-	1/1/1/1	-
4	EDO	A	503	-	-	1/1/1/1	-
4	EDO	C	509	-	-	1/1/1/1	-
4	EDO	D	519	-	-	1/1/1/1	-
4	EDO	C	504	-	-	1/1/1/1	-
4	EDO	A	511	-	-	0/1/1/1	-
4	EDO	D	509	-	-	0/1/1/1	-
4	EDO	C	503	-	-	1/1/1/1	-
4	EDO	B	503	-	-	1/1/1/1	-
4	EDO	D	504	-	-	1/1/1/1	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	EDO	A	512	-	-	1/1/1/1	-
4	EDO	B	510	-	-	1/1/1/1	-
4	EDO	D	514	-	-	1/1/1/1	-
4	EDO	D	503	-	-	1/1/1/1	-
4	EDO	D	505	-	-	1/1/1/1	-
4	EDO	D	520	-	-	1/1/1/1	-
4	EDO	D	510	-	-	1/1/1/1	-
7	PEG	D	512	-	-	1/4/4/4	-
4	EDO	C	508	-	-	1/1/1/1	-
4	EDO	B	505	-	-	0/1/1/1	-
4	EDO	D	513	-	-	0/1/1/1	-
5	E3Q	D	517	-	-	2/15/46/46	0/4/4/4
5	E3Q	C	507	-	-	6/15/46/46	0/4/4/4
4	EDO	A	506	-	-	1/1/1/1	-
4	EDO	B	504	-	-	1/1/1/1	-
7	PEG	D	511	-	-	3/4/4/4	-
5	E3Q	A	513	-	-	1/15/46/46	0/4/4/4
4	EDO	A	510	-	-	0/1/1/1	-
4	EDO	C	506	-	-	1/1/1/1	-
4	EDO	A	504	-	-	1/1/1/1	-
5	E3Q	B	509	-	-	3/15/46/46	0/4/4/4
4	EDO	A	508	-	-	1/1/1/1	-

All (12) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
6	B	508	EPE	C10-S	-10.67	1.62	1.77
6	C	505	EPE	C10-S	-8.93	1.65	1.77
6	D	516	EPE	C10-S	-7.34	1.67	1.77
6	B	511	EPE	C10-S	-6.96	1.67	1.77
5	D	517	E3Q	C24-N2	3.26	1.38	1.35
5	B	509	E3Q	C24-N2	3.07	1.38	1.35
5	B	509	E3Q	C17-N2	2.69	1.50	1.47
5	C	507	E3Q	C24-N2	2.62	1.37	1.35
5	A	513	E3Q	C24-N2	2.62	1.37	1.35
6	B	508	EPE	O2S-S	-2.33	1.38	1.45
5	D	517	E3Q	C17-N2	2.26	1.50	1.47
5	C	507	E3Q	C17-N2	2.08	1.49	1.47

All (36) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
6	D	516	EPE	O2S-S-C10	-14.08	85.45	106.73
6	D	516	EPE	O1S-S-C10	-12.81	87.38	106.73
6	D	516	EPE	O3S-S-C10	-9.50	87.43	106.00
5	A	513	E3Q	C16-N1-N2	-6.75	100.15	108.30
5	D	517	E3Q	C16-N1-N2	-6.38	100.60	108.30
5	B	509	E3Q	C16-N1-N2	-6.24	100.78	108.30
5	C	507	E3Q	C16-N1-N2	-6.20	100.83	108.30
5	C	507	E3Q	C11-C16-N1	-4.28	113.30	119.72
5	D	517	E3Q	C11-C16-N1	-4.18	113.44	119.72
5	B	509	E3Q	C11-C16-N1	-4.09	113.59	119.72
5	A	513	E3Q	C11-C16-N1	-3.76	114.08	119.72
5	A	513	E3Q	O2-C9-C14	3.59	123.17	115.75
6	B	511	EPE	O1S-S-C10	3.55	112.09	106.73
6	D	516	EPE	O3S-S-O1S	3.50	120.17	111.40
6	B	508	EPE	O1S-S-C10	3.32	111.75	106.73
6	D	516	EPE	O3S-S-O2S	3.22	119.45	111.40
5	A	513	E3Q	O2-C9-C10	-3.20	116.58	123.49
6	B	508	EPE	O2S-S-O1S	-3.02	103.99	113.82
5	A	513	E3Q	O1-C2-C7	2.69	119.97	115.96
6	B	508	EPE	O3S-S-C10	2.61	111.12	106.00
5	A	513	E3Q	O1-C2-C3	-2.61	119.91	124.30
5	C	507	E3Q	O1-C2-C7	2.47	119.63	115.96
6	C	505	EPE	O3S-S-O1S	-2.44	105.30	111.40
5	B	509	E3Q	O4-C24-N2	-2.43	124.96	127.01
5	C	507	E3Q	O4-C24-N2	-2.42	124.98	127.01
5	C	507	E3Q	O1-C2-C3	-2.40	120.25	124.30
5	C	507	E3Q	C17-N2-C24	-2.38	127.06	131.27
6	C	505	EPE	O1S-S-C10	2.29	110.18	106.73
6	B	511	EPE	O3S-S-C10	2.28	110.47	106.00
5	B	509	E3Q	C17-N2-C24	-2.23	127.32	131.27
6	C	505	EPE	O2S-S-C10	2.22	110.09	106.73
5	D	517	E3Q	O4-C24-C25	-2.21	124.80	126.93
5	D	517	E3Q	C17-N2-C24	-2.14	127.49	131.27
5	B	509	E3Q	O1-C2-C7	2.04	119.00	115.96
6	B	511	EPE	O3S-S-O1S	-2.03	106.31	111.40
6	C	505	EPE	C3-C2-N1	2.03	114.75	110.65

There are no chirality outliers.

All (59) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
6	B	508	EPE	C8-C7-N4-C3
6	B	511	EPE	C10-C9-N1-C6

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Mol	Chain	Res	Type	Atoms
6	B	511	EPE	C9-C10-S-O2S
6	B	511	EPE	C9-C10-S-O3S
6	C	505	EPE	C10-C9-N1-C2
6	C	505	EPE	C10-C9-N1-C6
6	C	505	EPE	C9-C10-S-O1S
5	C	507	E3Q	C3-C2-O1-C1
6	C	505	EPE	C9-C10-S-O3S
5	C	507	E3Q	C7-C2-O1-C1
6	D	516	EPE	N4-C7-C8-O8
7	D	511	PEG	O2-C3-C4-O4
4	A	507	EDO	O1-C1-C2-O2
4	A	508	EDO	O1-C1-C2-O2
4	C	503	EDO	O1-C1-C2-O2
4	D	507	EDO	O1-C1-C2-O2
4	D	519	EDO	O1-C1-C2-O2
5	A	513	E3Q	C7-C8-O2-C9
5	C	507	E3Q	C7-C8-O2-C9
5	B	509	E3Q	C7-C8-O2-C9
6	B	508	EPE	C9-C10-S-O3S
5	D	517	E3Q	C7-C8-O2-C9
7	D	511	PEG	O1-C1-C2-O2
6	B	508	EPE	N4-C7-C8-O8
4	B	510	EDO	O1-C1-C2-O2
4	C	504	EDO	O1-C1-C2-O2
4	D	505	EDO	O1-C1-C2-O2
4	D	510	EDO	O1-C1-C2-O2
4	D	515	EDO	O1-C1-C2-O2
4	A	504	EDO	O1-C1-C2-O2
7	D	511	PEG	C4-C3-O2-C2
6	B	508	EPE	C9-C10-S-O1S
6	B	508	EPE	C9-C10-S-O2S
6	B	511	EPE	C9-C10-S-O1S
6	C	505	EPE	C9-C10-S-O2S
6	B	508	EPE	S-C10-C9-N1
7	D	512	PEG	C1-C2-O2-C3
4	A	503	EDO	O1-C1-C2-O2
4	B	504	EDO	O1-C1-C2-O2
5	C	507	E3Q	C14-C9-O2-C8
5	B	509	E3Q	C14-C9-O2-C8
4	A	506	EDO	O1-C1-C2-O2
4	D	514	EDO	O1-C1-C2-O2
5	C	507	E3Q	C18-C17-N2-C24

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Mol	Chain	Res	Type	Atoms
4	A	505	EDO	O1-C1-C2-O2
4	B	503	EDO	O1-C1-C2-O2
4	B	506	EDO	O1-C1-C2-O2
5	D	517	E3Q	C14-C9-O2-C8
4	B	507	EDO	O1-C1-C2-O2
4	C	506	EDO	O1-C1-C2-O2
4	C	508	EDO	O1-C1-C2-O2
4	D	520	EDO	O1-C1-C2-O2
5	B	509	E3Q	C10-C9-O2-C8
4	A	512	EDO	O1-C1-C2-O2
4	D	503	EDO	O1-C1-C2-O2
4	D	504	EDO	O1-C1-C2-O2
4	C	509	EDO	O1-C1-C2-O2
4	D	508	EDO	O1-C1-C2-O2
5	C	507	E3Q	C10-C9-O2-C8

There are no ring outliers.

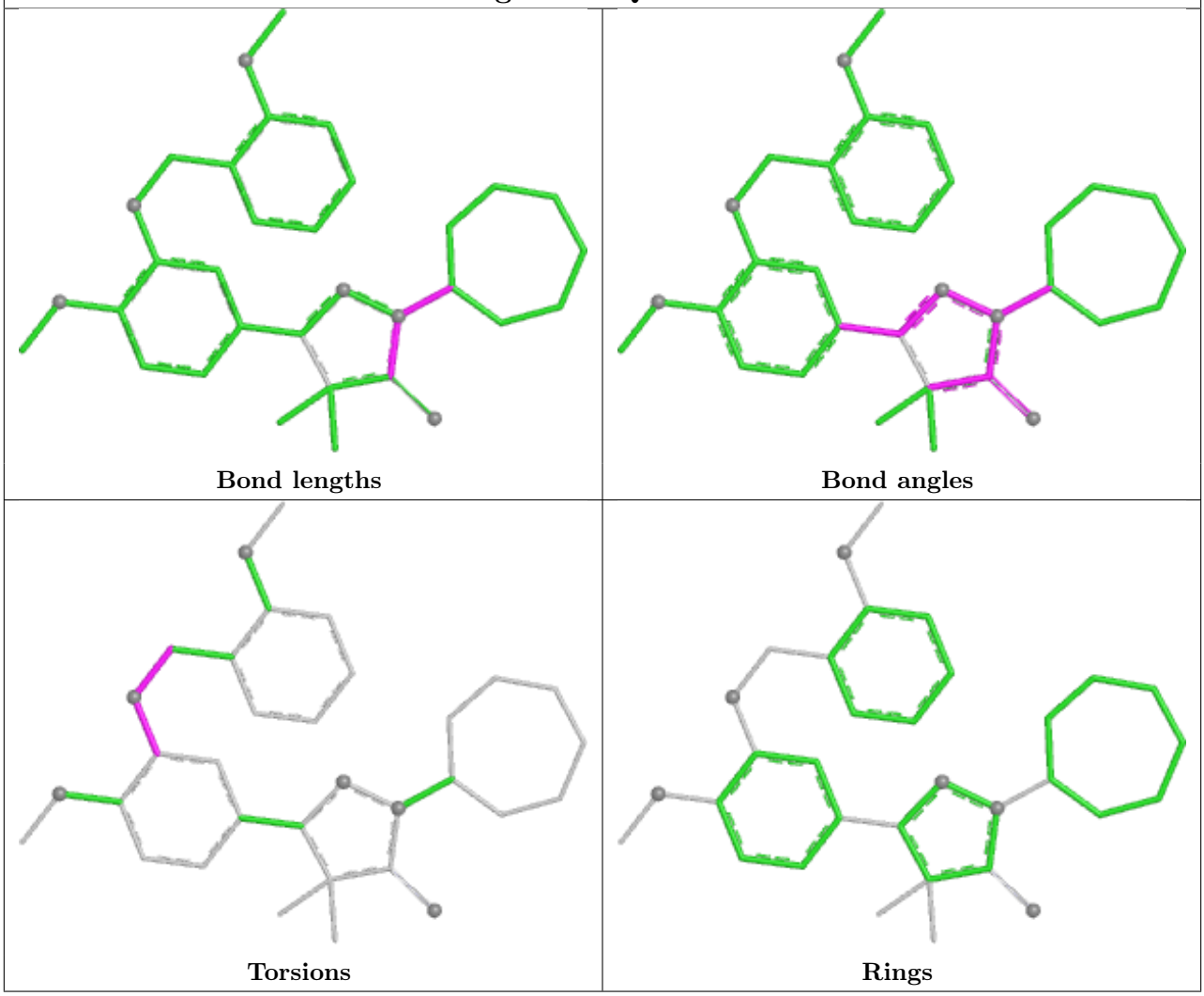
14 monomers are involved in 24 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	B	508	EPE	3	0
4	D	518	EDO	1	0
6	D	516	EPE	1	0
4	D	515	EDO	3	0
4	D	509	EDO	1	0
4	A	512	EDO	1	0
4	B	510	EDO	3	0
4	D	503	EDO	1	0
4	D	505	EDO	3	0
7	D	512	PEG	2	0
5	D	517	E3Q	1	0
7	D	511	PEG	1	0
5	A	513	E3Q	2	0
5	B	509	E3Q	1	0

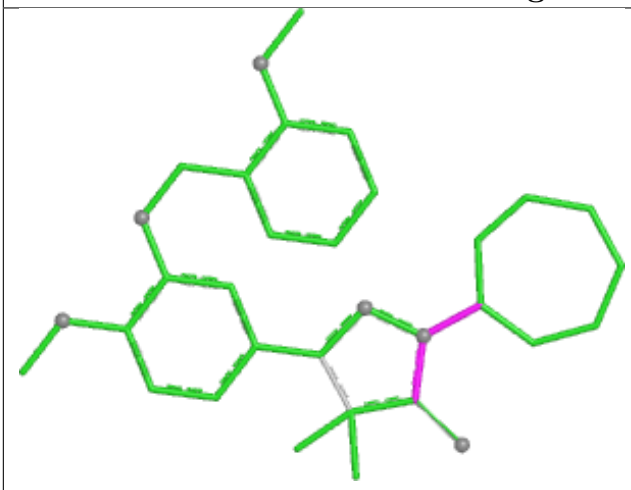
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 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.

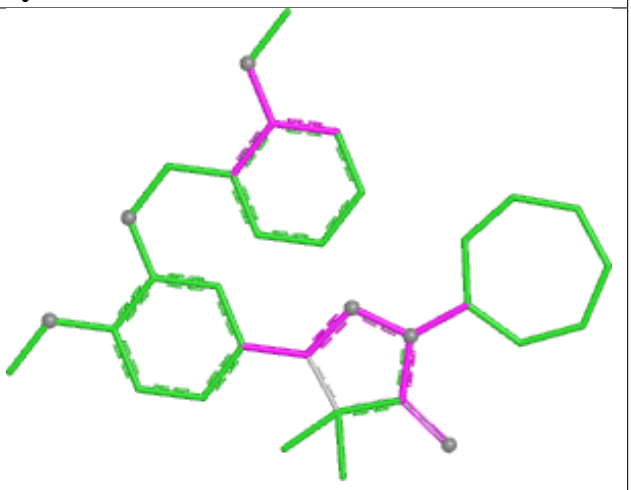
Ligand E3Q D 517



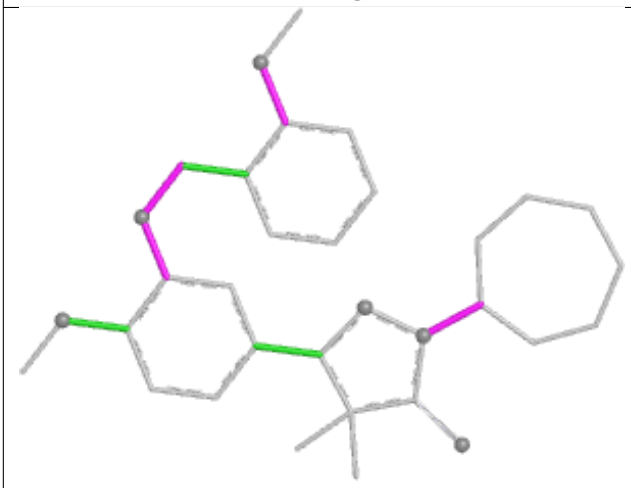
Ligand E3Q C 507



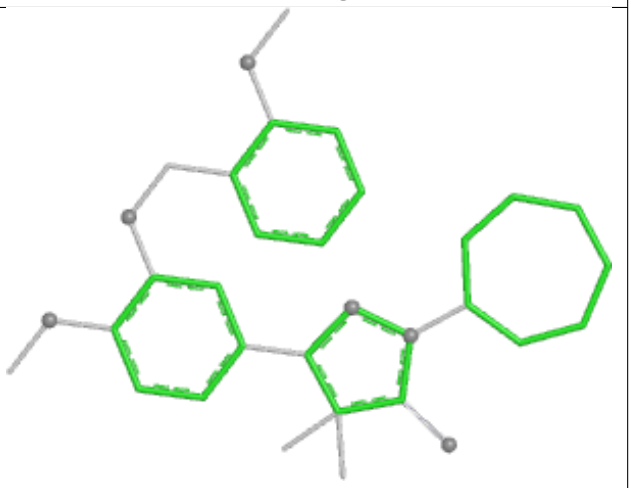
Bond lengths



Bond angles

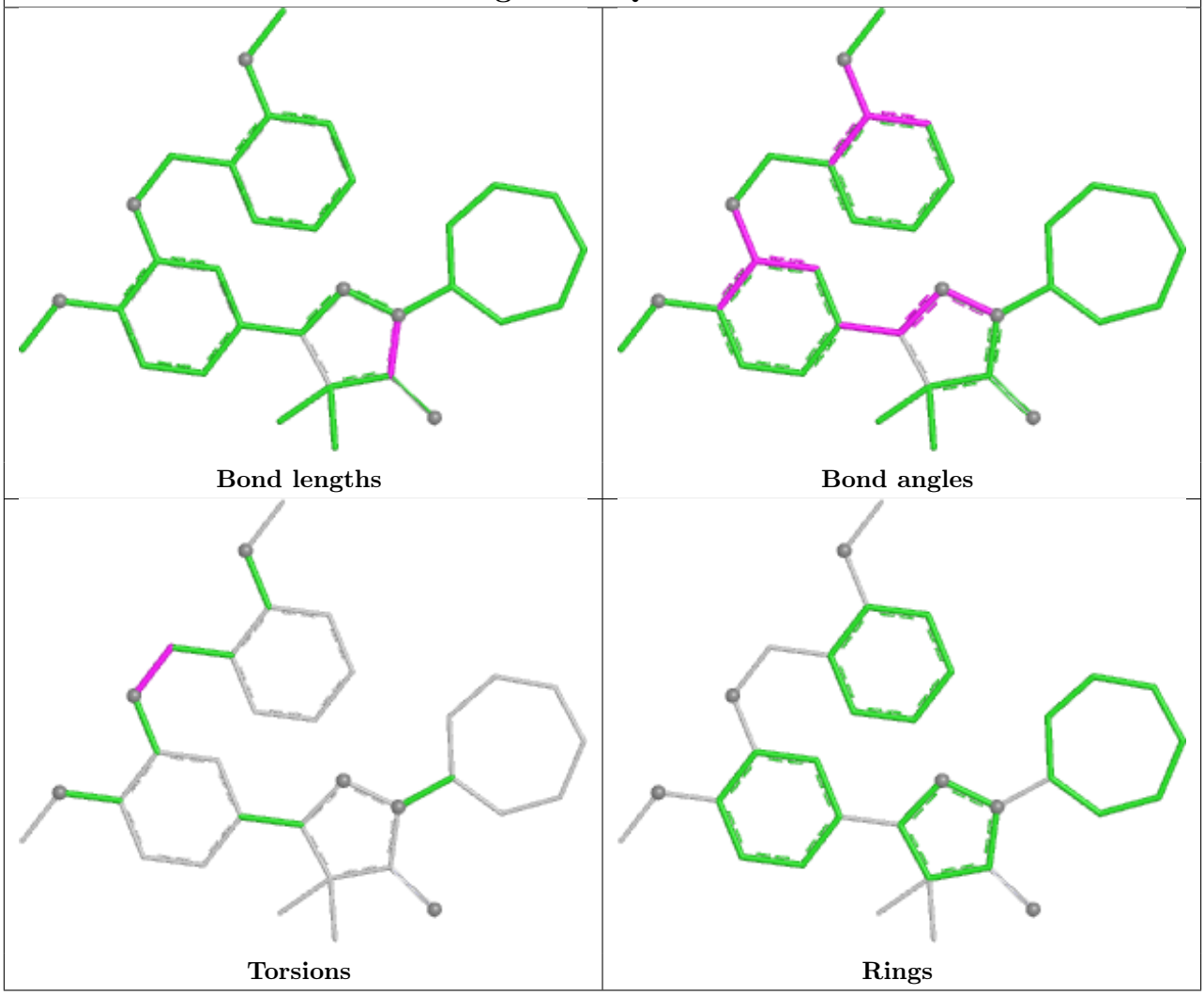


Torsions

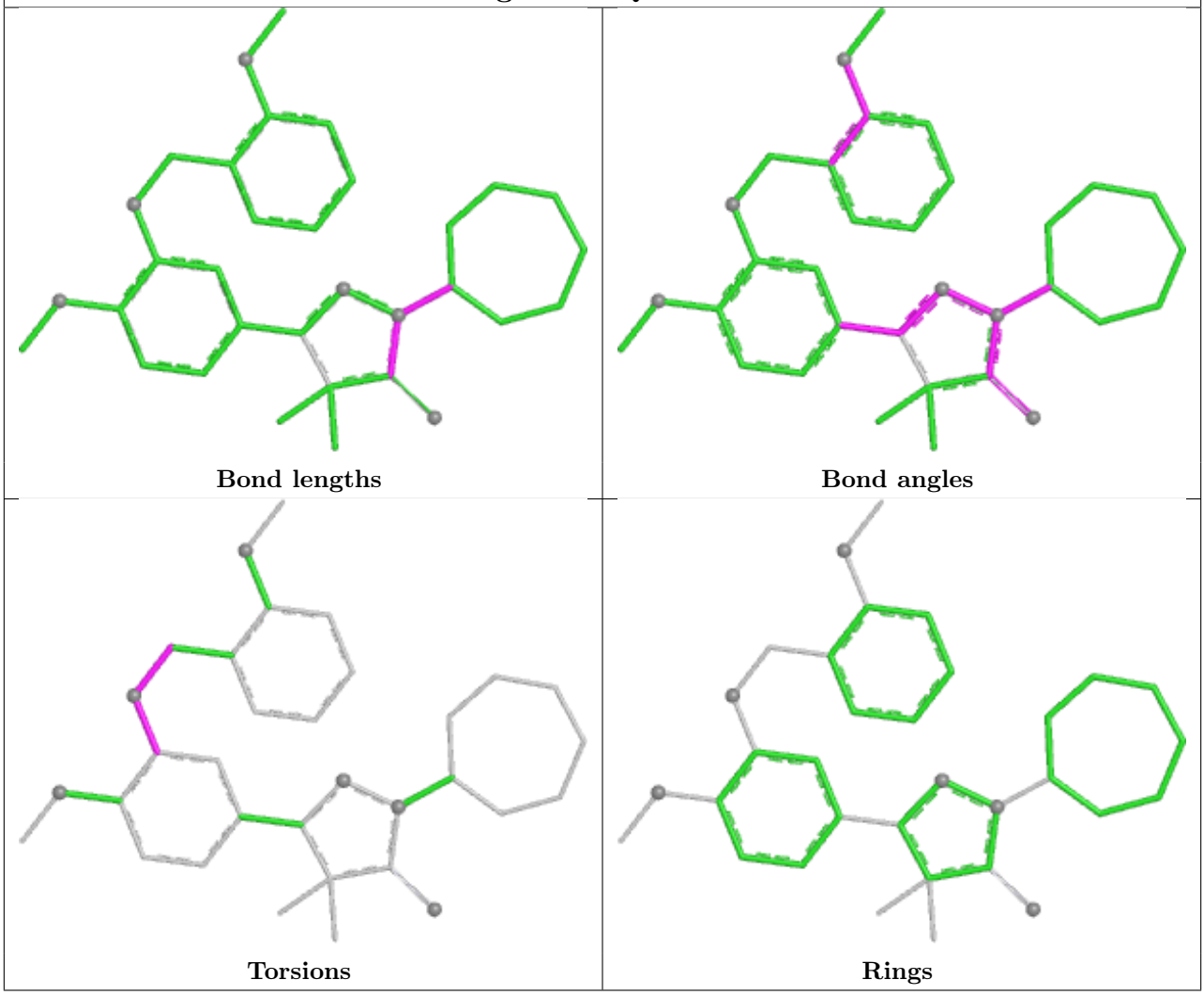


Rings

Ligand E3Q A 513



Ligand E3Q B 509



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 (i)

6.1 Protein, DNA and RNA chains (i)

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	329/364 (90%)	0.15	18 (5%)	25	24	19, 34, 71, 104	13 (3%)
1	B	324/364 (89%)	-0.08	1 (0%)	94	93	21, 37, 57, 88	0
1	C	323/364 (88%)	0.12	12 (3%)	41	41	21, 36, 71, 93	11 (3%)
1	D	324/364 (89%)	-0.08	2 (0%)	89	88	19, 28, 58, 98	0
All	All	1300/1456 (89%)	0.03	33 (2%)	57	56	19, 34, 65, 104	24 (1%)

All (33) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	358	CYS	11.9
1	A	357	MET	8.4
1	A	362	ASN	7.9
1	C	357	MET	6.7
1	A	352	MET	6.6
1	C	360	LYS	5.5
1	C	356	PRO	5.5
1	C	361	HIS	5.4
1	A	363	ALA	5.4
1	D	357	MET	5.3
1	A	351	GLY	4.7
1	A	356	PRO	4.6
1	A	360	LYS	4.4
1	A	354	ILE	4.0
1	A	350	ARG	3.9
1	C	359	ASP	3.8
1	C	411	PRO	3.6
1	A	358	CYS	3.6
1	C	354	ILE	3.5
1	A	340	PHE	3.4
1	C	355	SER	3.2

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Mol	Chain	Res	Type	RSRZ
1	A	349	GLU	3.0
1	A	411	PRO	2.8
1	C	363	ALA	2.8
1	A	353	GLU	2.6
1	A	355	SER	2.6
1	A	83	GLY	2.5
1	A	348	ARG	2.5
1	B	357	MET	2.4
1	C	362	ASN	2.4
1	D	354	ILE	2.2
1	A	294	SER	2.2
1	C	294	SER	2.1

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 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
4	EDO	C	508	4/4	0.69	0.18	53,53,53,62	0
4	EDO	D	506	4/4	0.71	0.25	74,77,78,80	0
5	E3Q	A	513	33/33	0.72	0.43	64,92,101,103	0
4	EDO	D	519	4/4	0.73	0.16	48,51,54,63	0
5	E3Q	B	509	33/33	0.79	0.23	39,59,81,89	0
4	EDO	C	504	4/4	0.80	0.23	56,56,57,61	0
4	EDO	A	508	4/4	0.83	0.17	58,63,63,70	0
4	EDO	D	507	4/4	0.83	0.24	34,42,45,46	0
4	EDO	D	513	4/4	0.83	0.20	55,56,60,61	0
4	EDO	B	504	4/4	0.84	0.24	53,54,54,61	0
4	EDO	D	515	4/4	0.85	0.18	41,44,44,46	0

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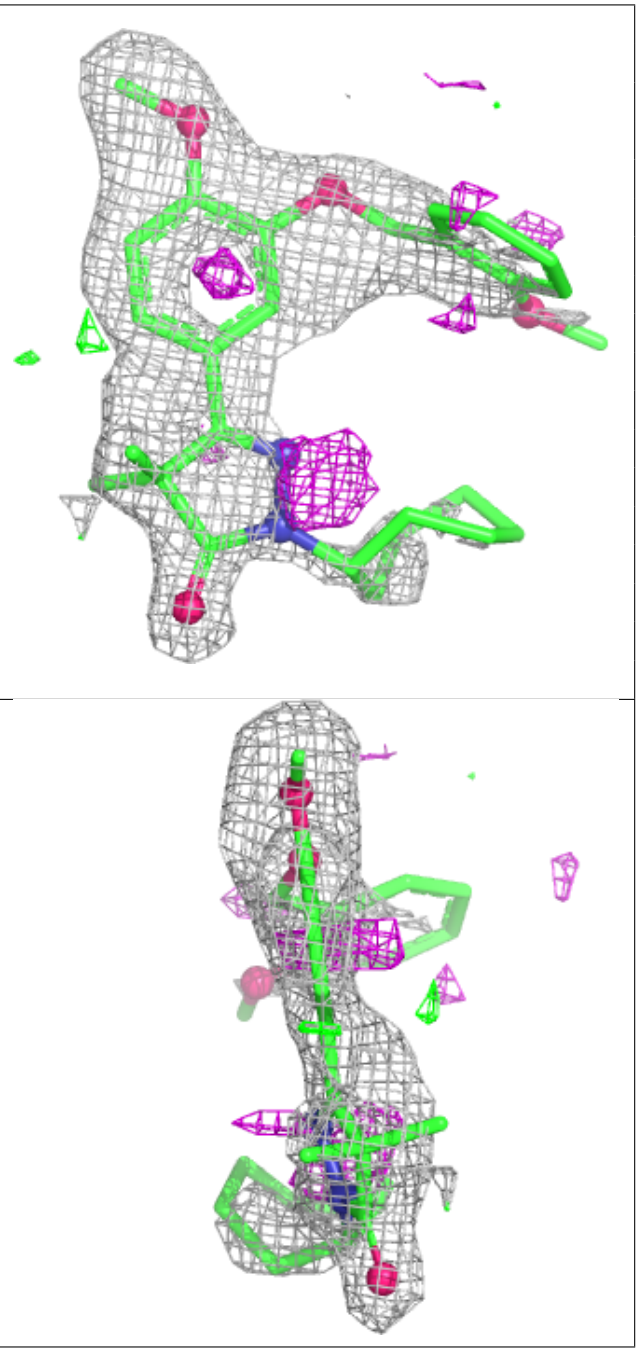
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	EDO	D	520	4/4	0.85	0.15	61,61,62,65	0
5	E3Q	C	507	33/33	0.85	0.26	52,64,83,85	0
5	E3Q	D	517	33/33	0.85	0.27	48,59,85,93	0
4	EDO	A	507	4/4	0.87	0.18	57,57,57,58	0
4	EDO	D	509	4/4	0.87	0.26	44,51,57,67	0
7	PEG	D	512	7/7	0.87	0.18	32,35,42,47	0
4	EDO	C	506	4/4	0.88	0.19	50,52,55,56	0
4	EDO	A	504	4/4	0.88	0.21	32,44,46,47	0
4	EDO	A	510	4/4	0.89	0.19	32,40,40,50	0
4	EDO	A	506	4/4	0.89	0.14	45,49,50,62	0
4	EDO	B	505	4/4	0.89	0.29	50,52,54,61	0
4	EDO	B	506	4/4	0.89	0.20	58,59,60,61	0
6	EPE	B	511	15/15	0.89	0.28	62,72,78,78	0
4	EDO	B	510	4/4	0.89	0.20	40,42,46,56	0
4	EDO	D	508	4/4	0.90	0.23	61,64,65,67	0
4	EDO	A	512	4/4	0.90	0.25	62,74,80,87	0
4	EDO	A	503	4/4	0.90	0.18	50,55,60,72	0
4	EDO	A	511	4/4	0.91	0.17	53,54,54,55	0
4	EDO	D	505	4/4	0.91	0.19	44,46,48,49	0
4	EDO	D	503	4/4	0.92	0.20	52,52,58,61	0
4	EDO	D	514	4/4	0.92	0.14	50,56,60,63	0
4	EDO	B	503	4/4	0.92	0.13	48,53,55,57	0
4	EDO	D	518	4/4	0.92	0.26	41,46,50,54	0
4	EDO	C	509	4/4	0.92	0.21	53,55,55,59	0
7	PEG	D	511	7/7	0.92	0.25	39,40,44,46	0
4	EDO	D	510	4/4	0.92	0.15	38,44,44,47	0
4	EDO	A	509	4/4	0.93	0.19	35,35,38,41	0
4	EDO	C	503	4/4	0.93	0.13	45,45,46,46	0
4	EDO	A	505	4/4	0.94	0.13	34,34,34,35	0
4	EDO	B	507	4/4	0.94	0.11	55,56,57,58	0
6	EPE	D	516	15/15	0.95	0.18	37,69,85,88	0
6	EPE	B	508	15/15	0.95	0.17	45,79,87,89	0
4	EDO	D	504	4/4	0.95	0.12	34,36,38,41	0
6	EPE	C	505	15/15	0.97	0.21	37,50,67,74	0
3	MG	A	502	1/1	0.97	0.13	19,19,19,19	0
3	MG	C	502	1/1	0.98	0.08	17,17,17,17	0
3	MG	B	502	1/1	0.98	0.09	17,17,17,17	0
2	ZN	B	501	1/1	0.99	0.09	31,31,31,31	0
2	ZN	C	501	1/1	1.00	0.09	31,31,31,31	0
2	ZN	D	501	1/1	1.00	0.11	26,26,26,26	0
3	MG	D	502	1/1	1.00	0.07	15,15,15,15	0
2	ZN	A	501	1/1	1.00	0.12	29,29,29,29	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.

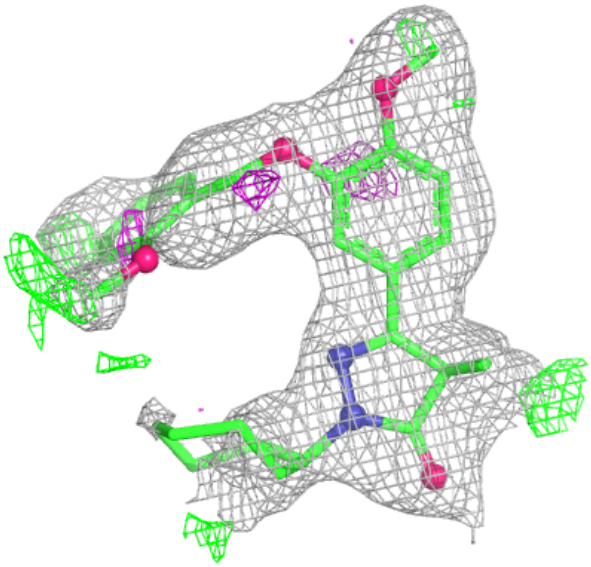
Electron density around E3Q A 513:

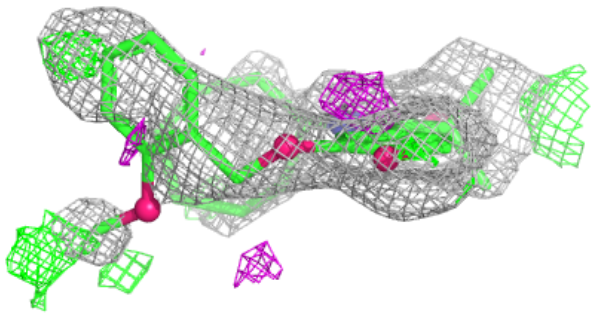
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)

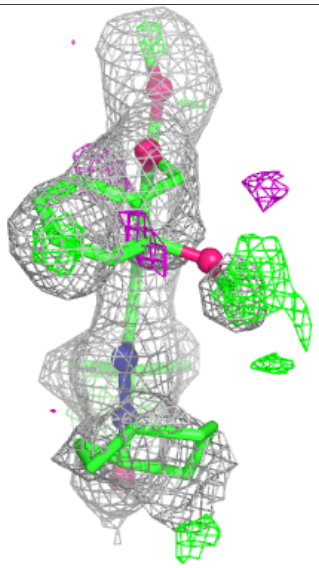


Electron density around E3Q B 509:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

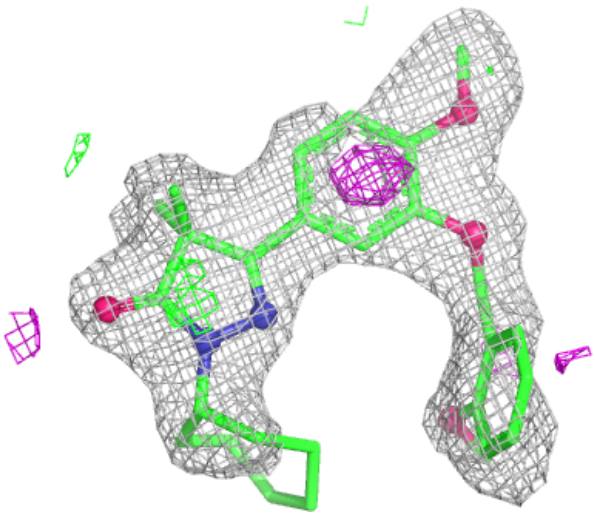


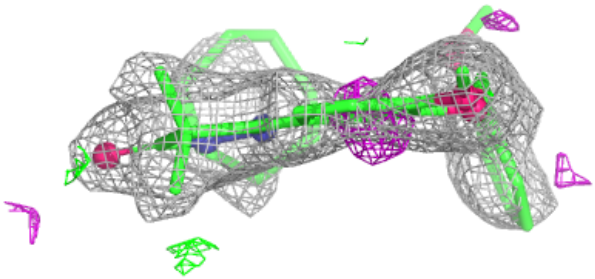


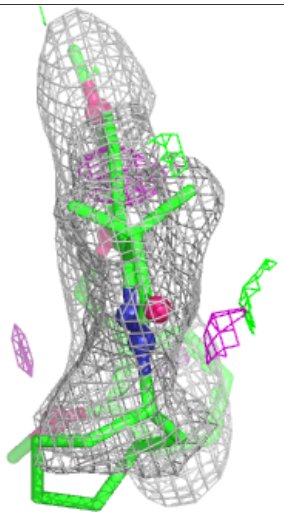


Electron density around E3Q C 507:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

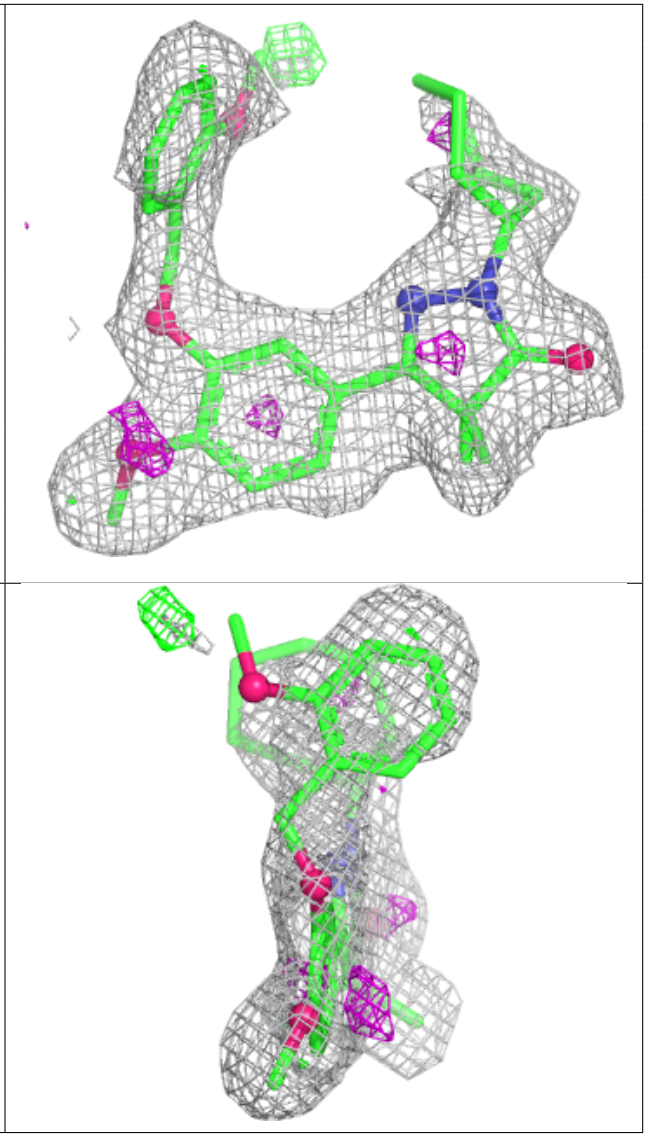






Electron density around E3Q D 517:

$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative) and green (positive)



6.5 Other polymers (i)

There are no such residues in this entry.